

A METHOD TO PERFORM FREE-VIBRATION ANALYSIS OF STRUCTURES

BY
HIRA LAL KAUL



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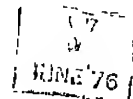
DEPARTMENT OF CIVIL ENGINEERING
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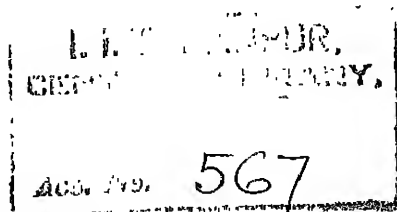
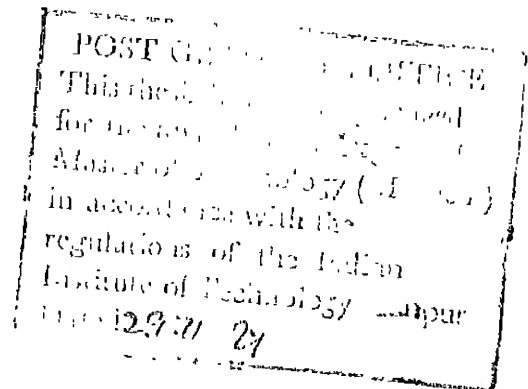
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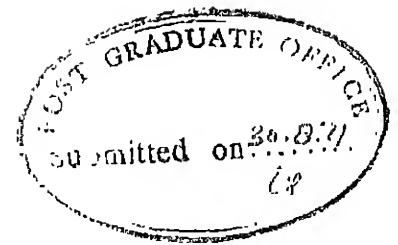


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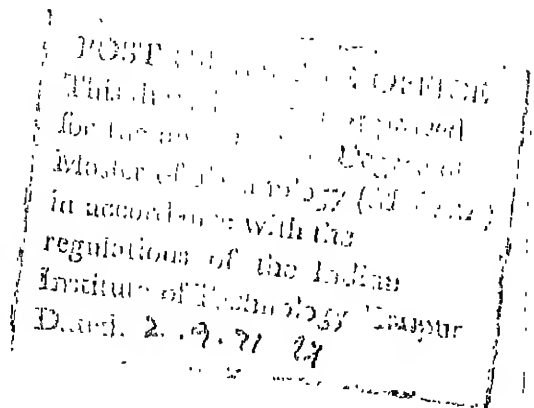
TO
MY PARENTS

CERTIFICATE



This is to certify that the thesis entitledd
'A Method to PERFORM FREE VIBRATION ANALYSIS OF STRUCTURES'
by Hira Lal Kaul is a record of work carried out under
my supervision and has not been submitted elsewhere for a
degree.

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SYNOPSIS
of the
Dissertation on
A METHOD TO PERFORM FREE VIBRATION
ANALYSIS OF STRUCTURES
A THESIS SUBMITTED
in partial fulfilment of the requirements
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The mode shapes, associated natural frequencies and nature and amount of damping govern the dynamic response of structures. This work presents a minimization method of extracting the eigenvalues and eigenvectors of a structural system having finite degrees of freedom. The structures considered are supposed to be linear and undamped. The Rayleigh quotient has been minimized successively in reduced subspaces to obtain the second and higher frequencies and the mode shapes associated with them. Conjugate Gradient Method of minimization has been used. Its modest demands on Computer storage space have been further supplemented

by performing the analysis which does not require the assembly of structural stiffness and mass matrices. The method seems to be particularly useful for large sized problems where only first few eigenvalues and eigenvectors are desired.

The study of free vibrations is usually an intermediate step to perform the forced vibration analysis. The method has been applied to simple structures like cantilever beam, portal frame and skew bridge and results obtained compare very well.

CHAPTER - I

INTRODUCTION

Performing a discrete element analysis of the structure, the matrix formulation of the general dynamic response problem for the undamped structure is:

$$[M] \ddot{\vec{Y}} + [K] \vec{Y} = \vec{F}(t) \quad \dots (1)$$

where the $n \times n$ matrices $[M]$ and $[K]$ are respectively the mass matrix and the stiffness matrix of the structure. These matrices are real and symmetric and their order n corresponds to the degrees of freedom to which an infinite degree of freedom continuous structure has been discretized. The $n \times 1$ vectors $\ddot{\vec{Y}}$, \vec{Y} and $\vec{F}(t)$ represent the absolute acceleration, absolute displacement and load respectively. The set of equations Eqn. (1), consist of a set of coupled second order ordinary differential equations. These are usually solved by uncoupling the equations through matrix transformation. The uncoupled equations resemble the equation of motion of a single degree of freedom system. The dynamic response is finally obtained by the linear superposition of the solution of these uncoupled equations. The matrix which performs uncoupling of the set of Eqn.(1)

requires the free vibration analysis of the structure which in turn means the solution to the set of equations:

$$[M] \ddot{\vec{Y}} + [K] \vec{Y} = 0 \quad \dots (2)$$

The above set of equations are reduced to an eigenvalue problem.

$$[K] \vec{X} = \lambda [M] \vec{X} \quad \dots (3)$$

where λ is the scalar quantity called eigenvalue which represents the square of the natural frequency of the system and \vec{X} is the $(n \times 1)$ eigenvector called the principal mode shape associated with the particular value λ .

This work presents a method to obtain the eigenvalues λ and associated eigenvectors \vec{X} . The method presented here is iterative and exploits the minimization technique and is expected to be specifically useful for determination of the first few eigenvalues and associated eigenvectors for large sized structures i.e. where n is large.

1.1 METHODS OF EIGENSOLUTION:

By eigensolution is meant the solution of the eigenproblem, Eqn. (3), which in turn means the determination of the scalar quantities λ and their associated

modeshapes \vec{X} .

Various methods of eigensolution known can be categorized as:

- (a) Transformation Methods
- (b) Iterative Methods.

Transformation methods require to convert the general eigenvalue problem:

$$[K]\vec{X} = [M]\vec{X} \text{ into a special form}$$

$$[A]\vec{X} = \lambda \vec{X}$$

and then involve a series of similarity transformations which operate on the matrix $[A]$. This requires the storage of the full $n \times n$ matrix $[A]$ in the computer. If n is large, it occupies sufficient computer space. Furthermore, all transformation methods like Givens, Householder, Jacobi etc., have been found to be efficient for the complete eigensolution i.e., determination of all the n eigenvalues and associated eigenvectors of the set of equations Eqn. (3), presuming that there exist n distinct eigenvalues.

However, a reasonably accurate dynamic response analysis of structures can be performed with the knowledge of only the first few eigenvalues and eigenvectors from the lowest end of the spectrum and hence the complete eigensolution is not required. Therefore, transformation methods are not very efficient for use in determining the dynamic response of large sized structures.

Iterative methods like Power Method, also require the initial conversion of general eigenvalue problem to a particular one and thus suffer with the same disadvantage as the transformation methods so far as computer space is concerned. Furthermore, these methods are not very stable and require periodic cleaning in order to obtain eigenvalues higher than the fundamental one.

The iterative method described and used in the present work has the advantage that it does not require any initial conversion of general eigenvalue problem to a particular form. Moreover, the method does not require the assembly of $(n \times n)$ structural stiffness and mass matrices. The entire algorithm uses the stiffness and mass matrices of the elements which compose the structure and their order is very small compared to n .

Lastly the method has been observed to be numerically stable and efficient for the determination

of partial eigensolution i.e., the first few natural frequencies and mode shapes of the vibrating structure.

Chapter 2 describes the formulation of the method for eigensolution. Chapter 3 describes the algorithm used to obtain the eigensolution. Chapter 4 gives the results for the class of structures, to which the method has been applied.

CHAPTER - II

FORMULATION OF THE METHOD OF EIGENSOLUTION*

An iterative method using the well known property of Rayleigh quotient is presented which can be applied directly to the eigenproblem:

$$[K] \vec{X} = \lambda [M] \vec{X}$$

It uses the property of the Rayleigh quotient:

$$R(\vec{X}) = \frac{\vec{X}^T [K] \vec{X}}{\vec{X}^T [M] \vec{X}}, \text{ that it equals the}$$

eigenvalue when the eigenvector is substituted into it and that it is stationary in the neighbourhood of an eigenvector. The Rayleigh quotient is minimized to obtain the lowest eigenvalue and the associated eigenvector. The minimization is done numerically using the conjugate gradient method. The intermediate eigenvalues and

* The method was developed and is described in detail in reference no. For completeness of presentation, it is briefly described here.

the associated eigenvectors are obtained by using a gradient projection scheme for constraining the minimization search to the subspace \mathcal{A} - orthogonal to the previously determined eigenvectors.

The advantage of the formulation

$$R(\vec{X}) = \frac{\vec{X}^T [\mathbf{K}] \vec{X}}{\vec{X}^T [\mathbf{M}] \vec{X}},$$

is that both the numerator and the denominator, as well as all of the other quantities needed for iteration procedure for all of the eigenvalue desired, can be computed without having the assembled $[\mathbf{K}]$ and $[\mathbf{M}]$ matrices at hand. Since, the numerator is twice the strain energy for a give \vec{X} (the displacement vector), and the denominator is twice the maximum kinetic energy of the structure, the summation of potential and kinetic energies of the individual elements of the discretized structural model gives us the numerator and the denominator.

2.1 FORMULATION OF THE PROBLEM:

The eigenproblem can be written as:

$$\left\{ [\mathbf{K}] - \lambda [\mathbf{M}] \right\} \vec{X} = 0$$

If \vec{X} is its solution, then $b \vec{X}$ is also a solution for any non-zero value of the scalar b , and thus

the eigenvector corresponding to any eigenvalue is arbitrary to the extent of a scalar multiplier. The Rayleigh quotient is not defined at the origin and thus care has to be taken while minimizing. The redundant degrees of freedom which prevents the determination of absolute magnitude of the components of the eigenvector, can be eliminated by an arbitrary normalization and the simplest normalization is to set any non-zero component of the eigenvector equal to one. The value of the Rayleigh quotient is bounded by the lowest and highest eigenvalues of the physical system.

Thus, the minimization of the Rayleigh quotient will yield the lowest eigenvalue. The minimization problem to yield the lowest eigenvalue can thus be stated as:

$$\begin{aligned} \text{Find } \vec{X} &= \vec{X}_1 \text{ such that} \\ R(\vec{X}_1) &= \frac{\vec{X}_1^T [\mathbf{K}] \vec{X}_1}{\vec{X}_1^T [\mathbf{M}] \vec{X}_1}, \end{aligned}$$

is minimum, subject to, $x_{q1} \equiv \vec{X}_1^T \vec{e}_q = 1$ where x_{q1} is the normalizing or reference component and \vec{e}_q is a vector with its q th component as one and zero elsewhere (i.e., \vec{e}_q is a unit co-ordinate vector for the q th co-ordinate).

Once the lowest eigenvalue $\lambda_1 \equiv R(\vec{X}_1)$ is known, the next higher or (second) eigenvalue and its associated eigenvector can be determined by posing a new minimization problem:

$$\begin{aligned} \text{Find } \vec{X} &= \vec{X}_2, \text{ such that} \\ R(\vec{X}_2) &= \frac{\vec{X}_2^T [\mathbf{K}] \vec{X}_2}{\vec{X}_2^T [\mathbf{M}] \vec{X}_2} \end{aligned}$$

is minimum, subject to

$$\vec{X}_2^T \vec{e}_r = 1$$

$$\text{and } \vec{X}_2^T [\mathbf{M}] \vec{X}_1 = 0.$$

In the subspace defined by the above constraints, the Rayleigh quotient takes on a unique minimum at the eigenvector associated with the second lowest eigenvalue. The second constraint represents the imposition of the M - orthogonality condition between \vec{X}_1 and \vec{X}_2 . Geometrically speaking, these constraints merely restrict the position of vector space in which the search for the second eigenvector is carried out and in this restricted subspace, R has a minimum corresponding to λ_2 .

Determination of the third and subsequent eigenvalues and the associated eigenvectors is accomplished by solving

a sequence of problems similar, as for the determination of 2nd eigenvalue. The only change is that each time one additional equation of constraint has to be imposed on the minimization problem to satisfy the condition that the eigenvector currently being sought is M - orthogonal to all of the previously determined eigenvectors. The problem of determining the l th eigenvalue ($2 \leq l \leq n$) can thus be written

$$\text{Find } \vec{X} = \vec{X}_l, \text{ such that}$$

$$R(\vec{X}_l) = \frac{\vec{X}_l^T [K] \vec{X}_l}{\vec{X}_l^T [M] \vec{X}_l}$$

$$\text{Subject to } \vec{X}_l^T \vec{e}_j = 1$$

$$\text{and } \vec{X}_l^T [M] \vec{X}_i = 0, \quad i = 1, 2, \dots, l-1$$

where \vec{X}_i , $i = 1, 2, \dots, l-1$ are assumed to be known when the l th eigenvector is being sought.

Denoting $[M] \vec{X}_i \equiv \vec{V}_i$, the above constraint equation can be written as

$$\vec{X}_l^T \vec{V}_i = 0, \quad i = 1, 2, \dots, l-1.$$

CHAPTER - III

MINIMIZATION - ALGORITHMS

The gradient methods are usually more powerful to minimize a function of n variables as they use the local information about the rate of change of function with respect to the changes in the variables. Hence one of the gradient methods is being used here for minimization of the Rayleigh quotient. The Rayleigh quotient as a function of the n variables $(X_1, X_2, \dots, X_n) \equiv \vec{X}$ is differentiable, and its gradient vector

$$\begin{aligned}\nabla R \equiv \vec{g} &= \frac{2 [\mathbf{K}] \vec{X}}{\vec{X}^T [\mathbf{M}] \vec{X}} - \frac{(\vec{X}^T [\mathbf{K}] \vec{X})}{(\vec{X}^T [\mathbf{M}] \vec{X})^2} \cdot 2 [\mathbf{M}] \vec{X} \\ &= \frac{2 ([\mathbf{K}] \vec{X} - R [\mathbf{M}] \vec{X})}{(\vec{X}^T [\mathbf{M}] \vec{X})}\end{aligned}$$

is easily computed.

The steepest descent method which is often faced with the convergence difficulties is modified for rapid convergence by conjugate gradient method. The Fletcher Reeves Conjugate gradient method has been used here, for reasons of its simplicity and its modest storage demands.

The algorithm is as follows:

3.1 FLETCHER AND REEVES METHOD:

Conjugate Gradients:

The convergence difficulties of the steepest descent (gradient) method can be greatly reduced by a very simple modification which converts it to the Conjugate Gradient method.

Here, the entire algorithm reduce to

$$\vec{x}_0 = \text{arbitrary}$$

$$\vec{G}_0 = \nabla F(\vec{x}_0)$$

$$\vec{S}_0 = -\vec{G}_0$$

$$\vec{x}_{i+1} = \vec{x}_i + \alpha_i^* \vec{S}_i$$

$$\vec{G}_{i+1} = \nabla F(\vec{x}_{i+1})$$

$$\beta_i = |\vec{G}_{i+1}|^2 / |\vec{G}_i|^2$$

$$\vec{S}_{i+1} = -\vec{G}_{i+1} + \beta_i \vec{S}_i$$

Clearly from this definition of \vec{S}_{i+1} , it is a linear combination of \vec{G}_{i+1} and $\vec{S}_0, \vec{S}_1, \dots, \vec{S}_i$

and hence, it is a linear combination of $\vec{G}_0, \vec{G}_1, \dots, \vec{G}_{i+1}$.

The conjugate gradient method is a rapidly convergent technique, suitable for use when the gradient of the function is readily computed. It generally far surpasses the steepest descent method except in unusual cases and with careful application, it is one of the most effective minimization techniques. In programming the method, attention is to be paid to the questions of the accuracy required in the determination of α^* , a strategy for restarting and a proper scaling.

There are more powerful methods, but they are generally impractical in large complicated problems.

The conjugate gradient method is the most efficient minimization technique as the problem size is increased, inspite of its weaker stability.

The problem of minimizing the Rayleigh quotient function to find the lowest eigenvalue is similar to an unconstrained minimization problem and the algorithm can be directly applied. The use of conjugate gradient method for finding the intermediate eigenvalue is possible only when the minimization of the Rayleigh quotient is restricted into a subspace of \vec{X} in which the constraints

$$\vec{X}_1^T \vec{e}_j = 1 \quad \text{and} \quad \vec{X}_1^T \vec{M} \vec{X}_i = 0, \quad i = 1, 2, \dots, n-1$$

is continuously satisfied. Thus, in order to insure that the search is carried out in the desired subspace of \vec{X} , it is necessary (1) to start the iteration with a point in that subspace and (2) to project the gradient vector \vec{g} , onto that subspace. Such a projection matrix is generated below:

3.2 PROJECTION MATRIX:

Let \vec{P} be a matrix which has the property that for any vector \vec{W} , the vector

$$\vec{W}_p \equiv \vec{P} \vec{W}$$

satisfies, $\vec{W}_p^T \vec{Z}_i = 0, \quad i = 1, 2, \dots, q$

where $\vec{Z}_i, i = 1, 2, \dots, q$ are q linearly independent vectors.

The above Eqn. can also be written in matrix form as

$$\begin{aligned} \begin{bmatrix} \vec{Z}_1 \\ \vdots \\ \vec{Z}_q \end{bmatrix}^T \vec{W}_p &= 0 \\ (q \times n) \quad (n \times 1) \end{aligned}$$

where $\begin{bmatrix} N \end{bmatrix}_{nxq} = \begin{bmatrix} \vec{z}_1, \vec{z}_2, \dots, \vec{z}_q \end{bmatrix}$

In otherwords, the matrix operator $\begin{bmatrix} P \end{bmatrix}$ eliminates from the vector \vec{W} the non-orthogonal components, thus giving the vector \vec{W}_p which is orthogonal to the subspace spanned by the vectors $\vec{z}_i, i = 1, 2, \dots, q$. This idea is differently expressed as:

$$\vec{W}_p = \vec{W} - \sum_{i=1}^q u_i \vec{z}_i, \text{ or in matrix form as}$$

$$\vec{W}_p = \vec{W} - \begin{bmatrix} N \end{bmatrix}_{(nxq)} \vec{u}_{(qx1)}$$

where the components of vector \vec{U} are $u_i, i = 1, 2, \dots, q$.

Pre-multiplying above Eqn. by $\begin{bmatrix} N \end{bmatrix}^T$, we obtain from

$$\begin{bmatrix} N \end{bmatrix}^T \vec{W}_p = 0$$

$$\begin{bmatrix} N \end{bmatrix}^T \vec{W}_p = \begin{bmatrix} N \end{bmatrix}^T \vec{W} - (\begin{bmatrix} N \end{bmatrix}^T \begin{bmatrix} N \end{bmatrix}) \vec{U} = 0$$

$$\therefore \vec{U} = (\begin{bmatrix} N \end{bmatrix}^T \begin{bmatrix} N \end{bmatrix})^{-1} \begin{bmatrix} N \end{bmatrix}^T \vec{W}$$

Thus, we obtain

$$\vec{W}_p = \vec{W} - \begin{bmatrix} N \end{bmatrix} (N^T N)^{-1} N^T \vec{W}$$

$$\vec{W}_b = \left\{ \begin{bmatrix} I \\ -[N] (N^T N)^{-1} N^T \end{bmatrix} \vec{W} \right.$$

$(N^T N)$ will be a $(q \times q)$ symmetric matrix, and is nonsingular since $[N]$ is a $(n \times q)$ matrix composed of q linearly independent columns.

Hence, its inverse exists.

The projection matrix, by comparison is obtained as

$$[P] = \left\{ [I] - [N] ([N]^T [N])^{-1} [N]^T \right\}$$

For determining the l th eigensolution, a projection matrix

$$[P_1] = \left\{ \begin{bmatrix} I \\ -[N_1] ([N_1]^T [N_1])^{-1} [N_1]^T \end{bmatrix} \right\}$$

$(n \times 1) \quad (1 \times 1) \quad (1 \times n)$

where $[N_1] = [\vec{e}_j, \vec{v}_1, \vec{v}_2, \dots, \vec{v}_{l-1}]$

will project the gradient vector \vec{g} , on to the subspace of constraints already defined. The column vectors of the matrix $[N_1]$ are linearly independent. Thus

$$\vec{g}_p = [P_1] \vec{g}$$

One thing that warrants attention here is that, since we use \vec{g}_p instead of \vec{g} , will the conjugate gradient method actually converge? The answer is

affirmative and it can be shown by constructing the Lagrange function for the unconstrained minimization problem.

STEP BY STEP PROCEDURE:

3.3 (a) Choice of a Starting Point:

A the iterative methods should have a good starting point, otherwise time is unnecessarily wasted inside the minimization procedure. A starting vector as a unit vector \vec{e}_i (the j th element of \vec{e}_i is δ_{ij}) lying along the co-ordinate axes, gives an awfully distorted mode shape in the dynamics problem. A set of random vectors proved to be superior contrary to the expectation and in this work they were used as starting vectors. The starting point for the search of the lowest eigenvalue needs to satisfy only one constraint $\vec{x}_{q1} \equiv \vec{X}_1^T \vec{e}_q = 1$ which is trivially satisfied by dividing through out by the q th component. The minimization algorithm using conjugate gradient method generate a sequence of vector which, in the limit, tend directionally to the minimum eigenvalue on the search space. Thus the q th component of the vectors so generated have to be maintained as unity, so as to satisfy the constraint Eqn. $\vec{x}_{q1} = \vec{X}_1^T \vec{e}_q = 1$ continuously in the space. This is achieved by setting

the q th component of the gradient vector at the particular point equal to zero. In other words, no 'move' is made in the q th direction of the search space. The starting point for the search of second eigenvalue has to satisfy an additional constraint Eqn. $\tilde{\mathbf{X}}_2^T [\mathbf{M}] \tilde{\mathbf{X}}_1 = 0$, and this is achieved by Schmidt orthogonalization.

Let $\tilde{\mathbf{X}}(0)$ be some initial estimate.

Therefore, $\tilde{\mathbf{X}}_2(0) = \tilde{\mathbf{X}}_2(0) - (\tilde{\mathbf{X}}_2(0)^T \tilde{\mathbf{V}}_1) \tilde{\mathbf{V}}_1$

satisfies the above constraint if $\tilde{\mathbf{V}}_1^T \tilde{\mathbf{V}}_1 = 1$.

Since $\tilde{\mathbf{V}}_1 \equiv [\mathbf{M}] \tilde{\mathbf{X}}_1$, where $\tilde{\mathbf{X}}_1$ is the eigenvector corresponding to the first (lowest) eigenvalue.

The starting point for the search of the subsequent eigenvalues, is obtained by passing the initial estimate through a projection matrix, where the projection has for

$$[\mathbf{M}_1], \text{ the following}$$

$$[\tilde{\mathbf{M}}_1] \equiv [\tilde{\mathbf{V}}_1, \tilde{\mathbf{V}}_2, \dots, \tilde{\mathbf{V}}_{1-1}]$$

($n \times l-1$)

3.4 (b) FUNCTION EVALUATION:

It is necessary to have an efficient routine for function evaluation, in order to avoid the time which would be otherwise wasted inside the minimization procedure.

Now, $R(\vec{X})$ can be written as

$$R(\vec{X}) = \frac{\sum_{i=1}^r \vec{Y}_i^T [k_i] \vec{Y}_i}{\sum_{i=1}^r \vec{Y}_i^T [m_i] \vec{Y}_i}$$

where r is the number of discrete elements, $[k_i]$ and $[m_i]$ are, respectively the stiffness and mass matrices of the i th element of the discretized structure and \vec{Y}_i is the displacement vector of the i th element corresponding to the generalized displacement vector \vec{X} .

The size of element stiffness and mass matrices are relatively much smaller than the size of the assembled stiffness and mass matrices of large complex structure. The stiffness and mass matrices of various kinds of one dimensional structural elements are given in Appendix - II.

3.5 (c) Gradient Evaluation:

The minimization algorithm require the evaluation of gradient vector at each cycle of the iteration. The gradient vector, \vec{g} , of the Rayleigh quotient function is given earlier. For determining the lowest eigensolution, the component of the gradient vector corresponding to the normalizing component of the eigenvector is set equal

to zero, in order to continuously satisfy the constraint imposed due to the normalization of the eigenvector. For the intermediate eigensolution, the projection matrix $[P_1]$ is used to project this gradient vector, \vec{g} , on to the proper subspace of search, μ - orthogonal to the previously determined eigenvector. The component of the gradient vector, \vec{g}_p , corresponding to the normalizing component of the eigenvector turns out to be zero automatically, but a small number generally appears because of roundoff errors and this is removed by setting that component equal to zero.

3.6 (d) Evaluation of Step Length:

Once a direction of move \vec{S}_i has been decided, we must determine α_i^* so that the function is minimized in that direction. Thus the problem of finding the step length is essentially the linear search problem which requires the determination of the α_i^* along \vec{S}_i through \vec{X}_i at which the value of the Rayleigh quotient function

$$R(\vec{X}_i + \alpha_i \vec{S}_i) = \frac{(\vec{X}_i + \alpha_i \vec{S}_i)^T [K] (\vec{X}_i + \alpha_i \vec{S}_i)}{(\vec{X}_i + \alpha_i \vec{S}_i)^T [M] (\vec{X}_i + \alpha_i \vec{S}_i)}$$

is a minimum, i.e.,

$$\left. \frac{dR/d\alpha_i}{\alpha_i = \alpha_i^*} \right|_{\alpha_i = \alpha_i^*} = R'(\alpha_i^*) = 0$$

In the general problem, no expression is available to determine α_i^* , so an interpolation approach is adopted. However, in the particular problem of the Rayleigh quotient, an explicit expression in α_i can be generated, which has the form

$$u\alpha_i^2 + v\alpha_i + w = 0$$

$$\text{where } u = (\vec{S}_i^T [K] \vec{S}_0) (\vec{X}_i^T [M] \vec{S}_i) \\ - (\vec{X}_i^T [K] \vec{S}_i) (\vec{S}_i^T [M] \vec{S}_i)$$

$$v = (\vec{S}_i^T [K] \vec{S}_i) (\vec{X}_i^T [M] \vec{X}_i) - (\vec{X}_i^T [K] \vec{S}_i) (\vec{S}_i^T [M] \vec{S}_i)$$

$$w = (\vec{X}_i^T [K] \vec{S}_i) (\vec{X}_i^T [M] \vec{X}_i) - (\vec{X}_i^T [K] \vec{X}_i) (\vec{X}_i^T [M] \vec{S}_i)$$

The two roots correspond to the maximal and minimal points of the Rayleigh quotient in the direction \vec{S}_i through \vec{X}_i . The minimal function value corresponds to α_i^* .

CHAPTER - IV

APPLICATION - AND - RESULTS

In all the examples presented in this Chapter, the stiffness and mass matrices are taken from Appendix - II. The distributed mass of the system was used to evaluate the components of the mass matrix (the 'consistent' mass matrix). IBM 7044 digital computer was used to obtain the numerical results and random vectors were taken as starting points for the minimization algorithm to obtain the eigensolution.

4.1 EXAMPLE (1):

As a simple application, a slender solid cantilever rod shown in Fig. (1) is considered. The first three eigenvalues and eigenvectors of this simple structure were obtained by the successive minimization of the Rayleigh quotient in the appropriate subspace.

The cantilever rod shown in Fig.(1) is divided into three elements and each of them is modelled as a general planar beam element having six degrees of freedom. The cantilever Fig.(1) thus has a total of nine degrees of freedom. The first three eigenvalues and their associated mode shapes are given in Fig. 2(a) to 2(c).

The Table (1) / compares the computed results with actual theoretical results obtained from closed form analytical expression.

4.2 Another illustrative example considered here is the portal frame shown in Fig. 3. Each member was modelled with a single beam element. Thus the structure has four degrees of freedom. A complete eigensolution was obtained by successive minimization of the Rayleigh quotient. The time taken was 24 seconds.

The computed frequencies are given in Table 2.

4.3 Finally a skew bridge structure with 12 degrees of freedom shown in Fig. 5, has been analyzed to obtain the first four eigenvalues and the associated eigenvectors by successive minimization of R.Q. The analysis is carried out for a grillage with two longitudinals, fixed at the extreme ends and their cross girders are assumed to be simply supported over them. The torsional rigidity of the members is neglected.

The computed values of frequencies obtained are compared with calculated values of the frequencies and the experimental values of frequencies as given in Ref. (11),

and are indicated in Table 3.

The first mode shape is plotted and is shown in Fig. 6.

The total time taken to compute the first four eigenvalues and the associated eigenvectors is four minutes approximately.

TABLE 1

Frequency No.	Actual Values in cs./sec.	Computed values in cs/sec.
1	5.16	5.099
2	31.599	31.799
3	88.38	88.99

TABLE 2

COMPUTED FREQUENCIES IN CS. /SEC. FOR PORTAL
FRAME ARE:

1. 42.99
2. 241.25
3. 833.23
4. 1559.72

TABLE 3

26

Frequency No.	Calculated value of f. in cs/sec.	Experimental value of f in cs/sec	Computed value in cs/sec.
1	11.56	11.222	11.605
2	12.72	12.53	12.80
3	27.82	27.98	28.30
4	35.44	34.21	36.62

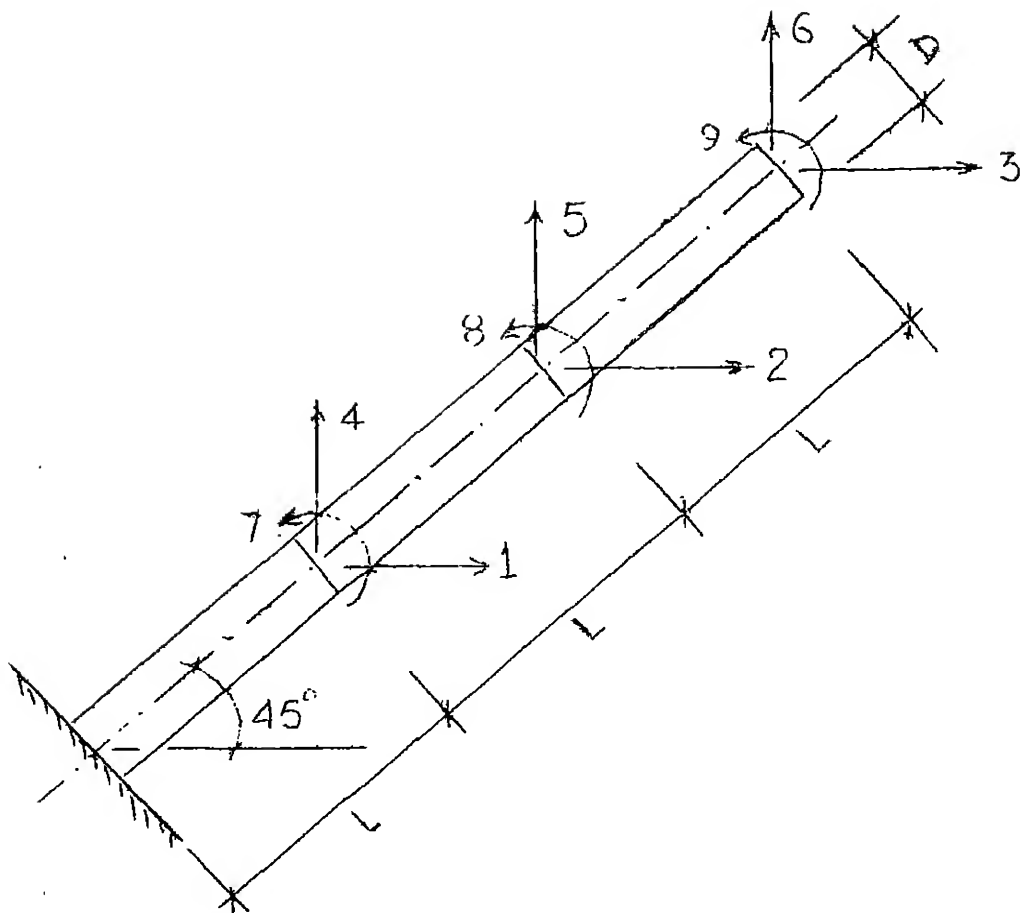


FIG. 1
SLENDER CANTILEVER ROD (9D.O.F.)

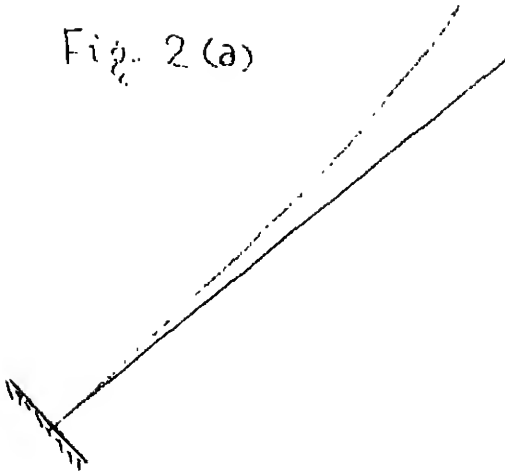
$$L = 25.0 \text{ in.}$$

$$D = 1.0 \text{ in.}$$

$$E = 30.0 \times 10^6 \text{ lbs./in}^2$$

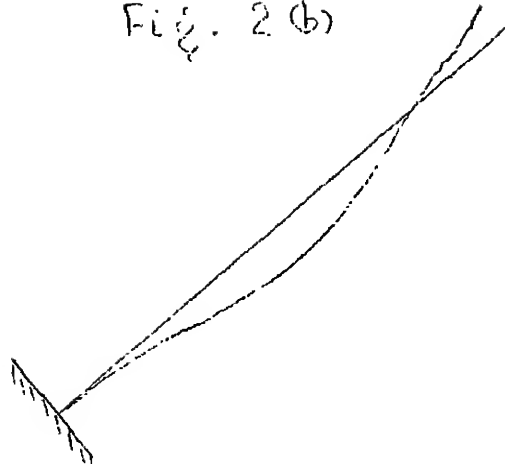
$$w = 0.28 \text{ lbs/in}^2.$$

Fig. 2 (a)



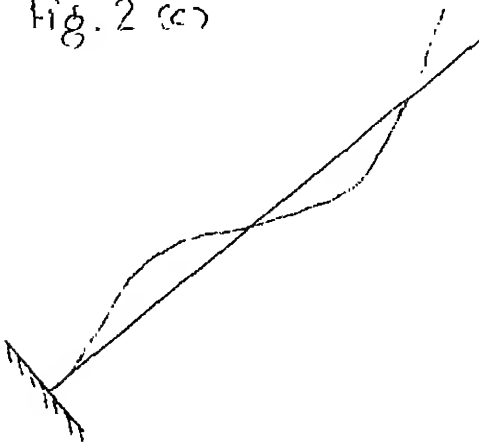
FIRST MODAL SHAPE

Fig. 2 (b)



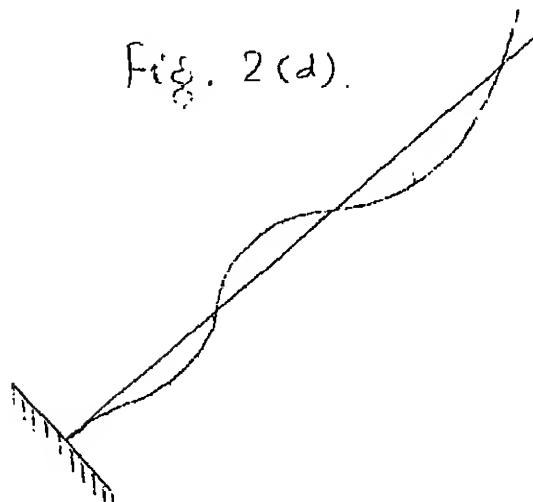
SECOND MODAL SHAPE

Fig. 2 (c)



THIRD MODAL SHAPE

Fig. 2 (d)



FOURTH MODAL SHAPE

CANTILEVER MODE SHAPES

FIG. 2

(Refer Fig 1 - a slender Cantilever Rod inclined at 45°)

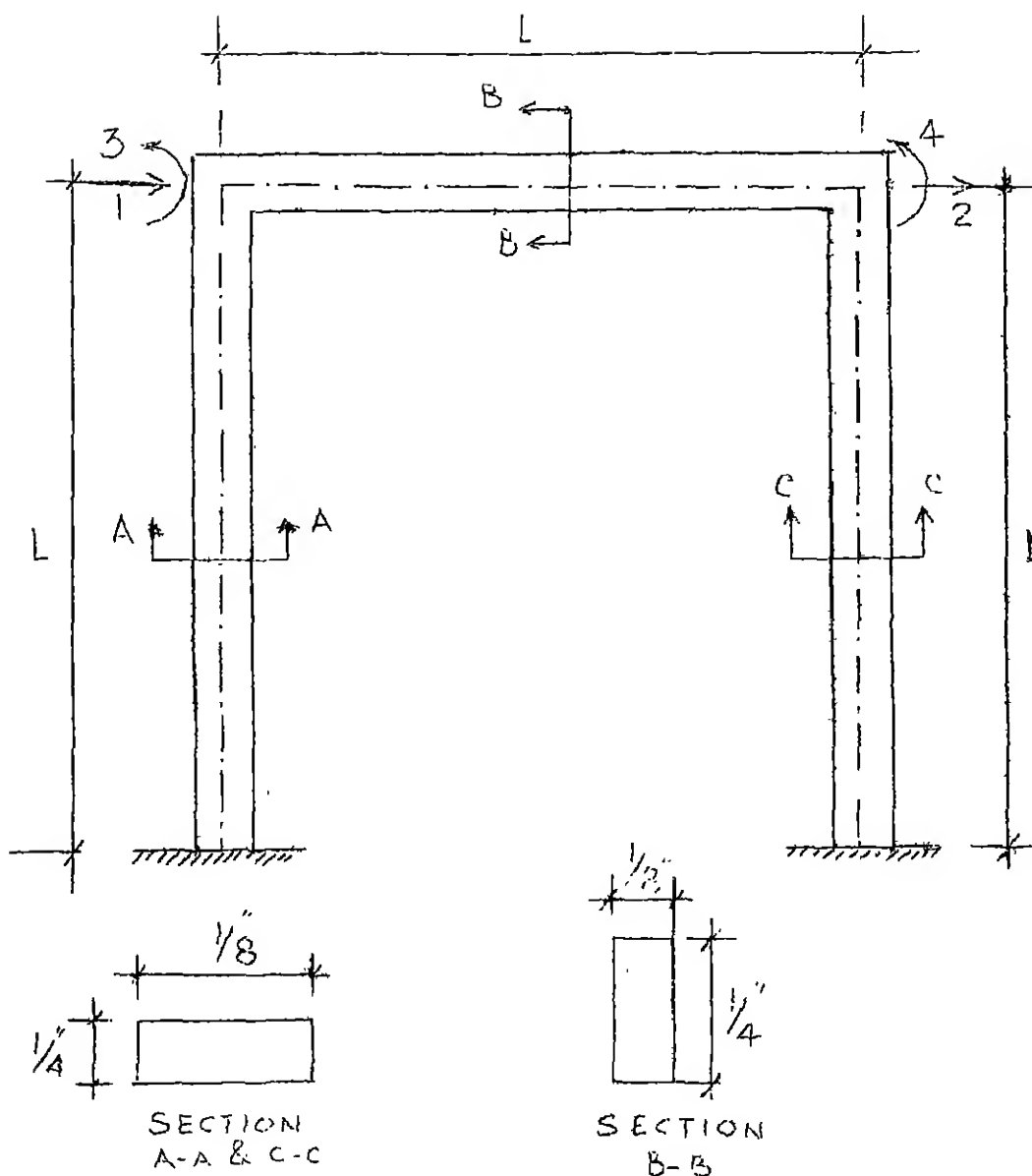


FIG. 3

PORTAL FRAME ; FOUR DEGREES OF FREEDOM SYSTEM.

$$L = 25.4 \text{ cm (10 inches)}$$

$$E = 2.1 \times 10^6 \text{ kgs/cm}^2 \text{ (} 30.0 \times 10^6 \text{ lbs/in}^2 \text{)}$$

$$= 7.81 \text{ gm/cm}^3 \text{ (0.28 lbs/inch}^3 \text{)}$$

DIFFERENT MODE SHAPES
OF A PORTAL FRAME.

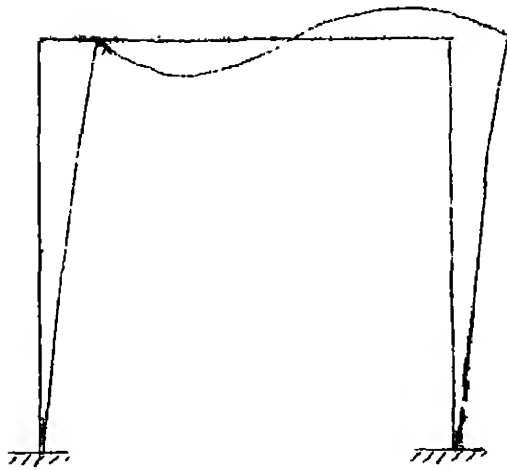


FIG. 4 (a)
FIRST MODAL SHAPE

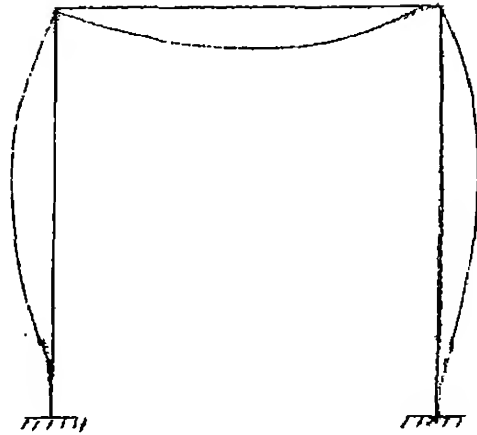


FIG. 4 (b)
SECOND MODAL SHAPE

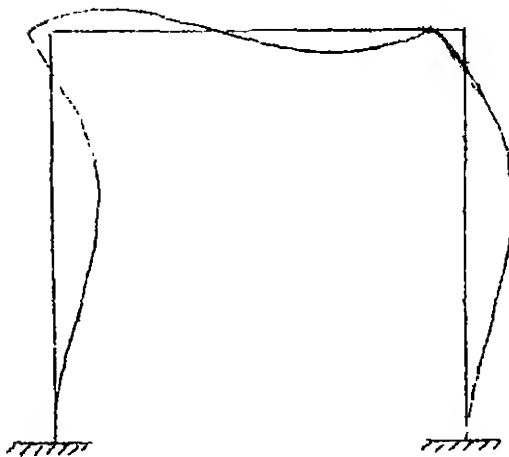


FIG. 4 (c)
THIRD MODAL SHAPE

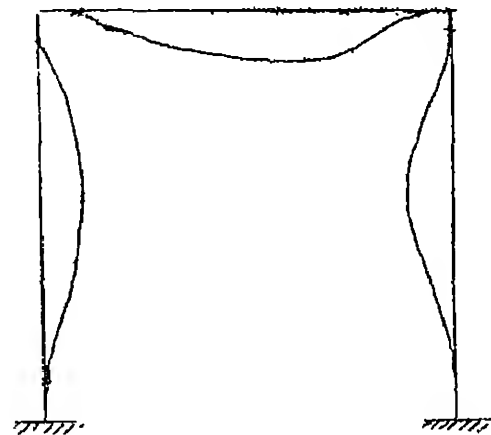


FIG. 4 (d)
FOURTH MODAL SHAPE

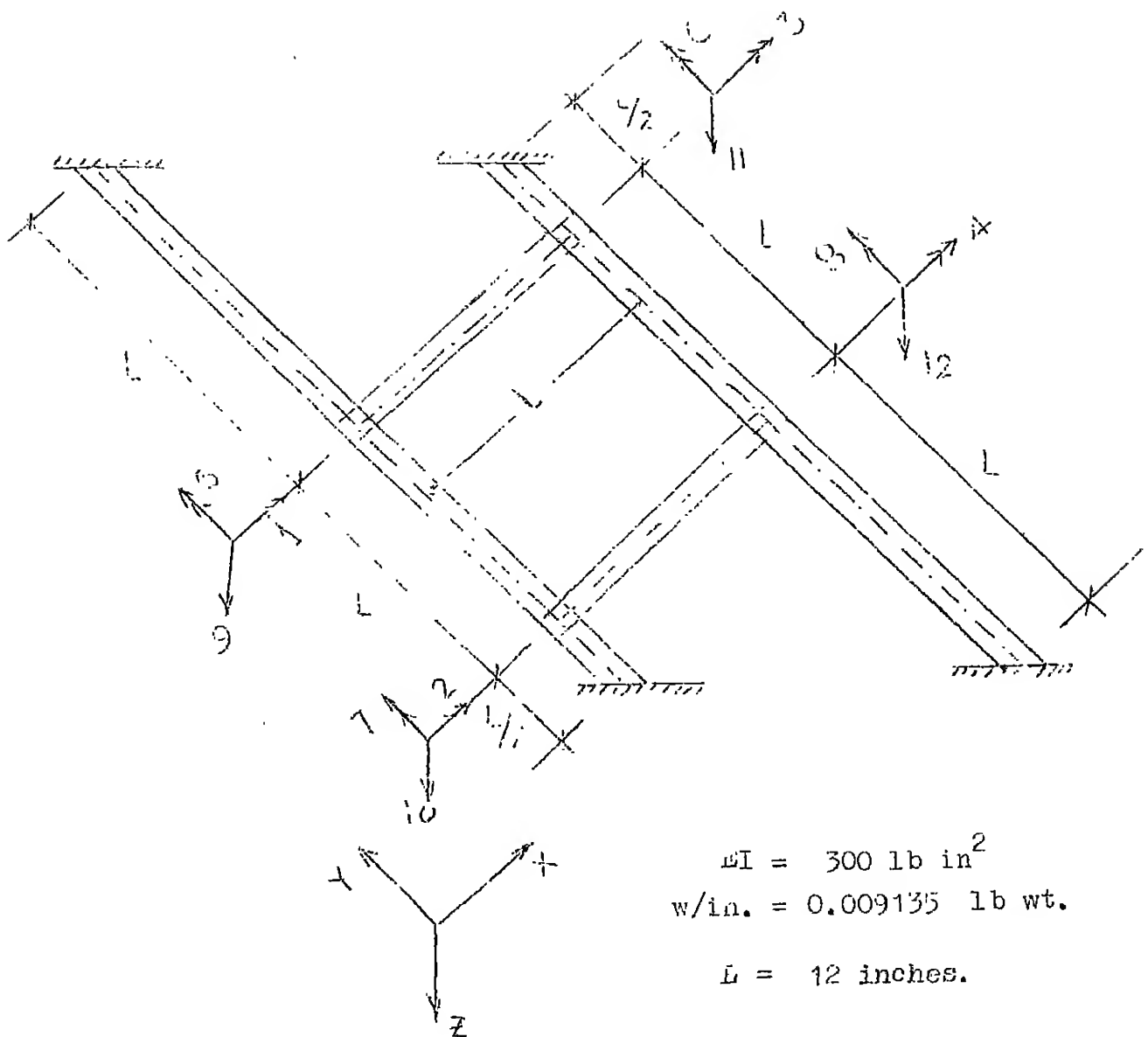
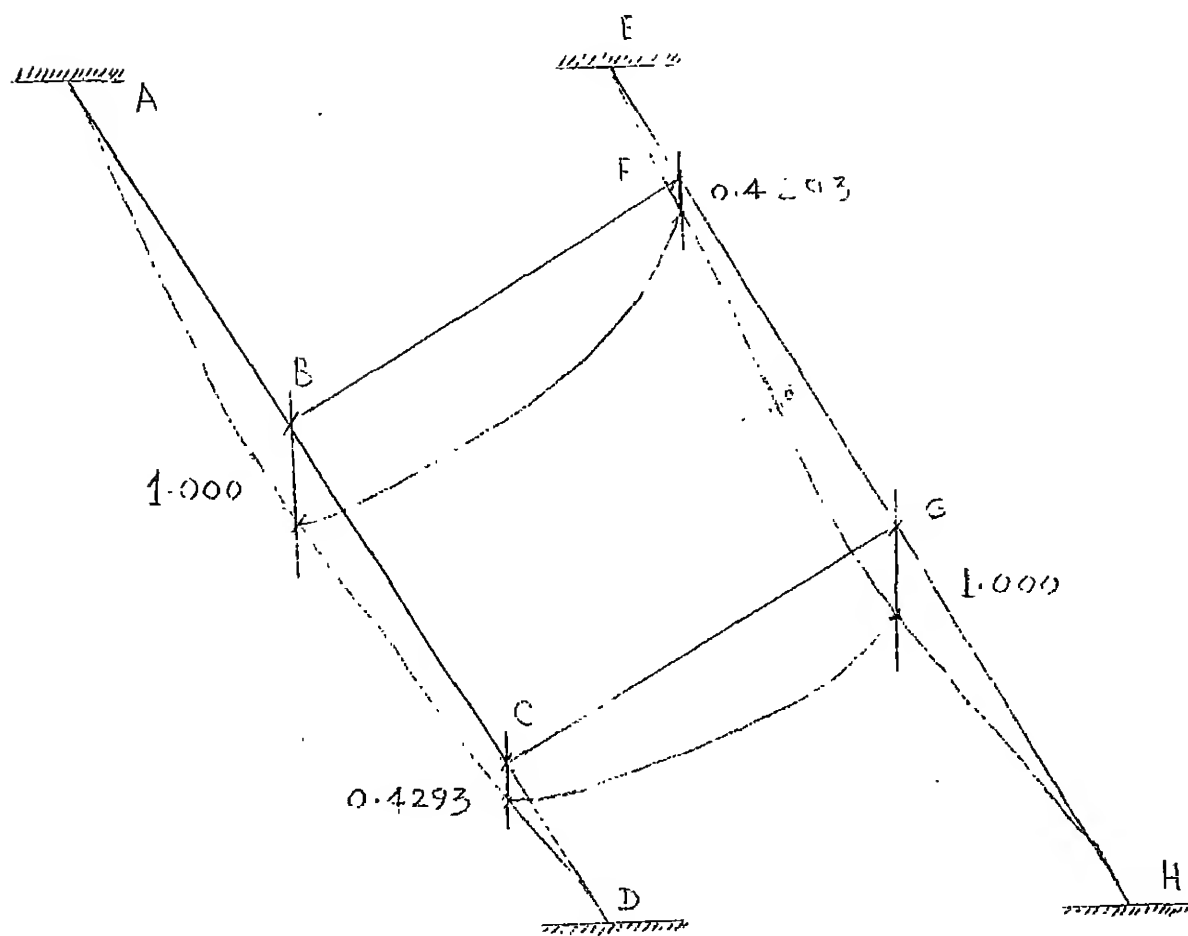


FIG. 5

SKW BRIDGE; 12 DEGREES O. FREEDOM

MODAL SHAPE OF A SKEW BRIDGE



First Mode Shape of a Skew Bridge
shown in Fig. 5.

FIG. 6

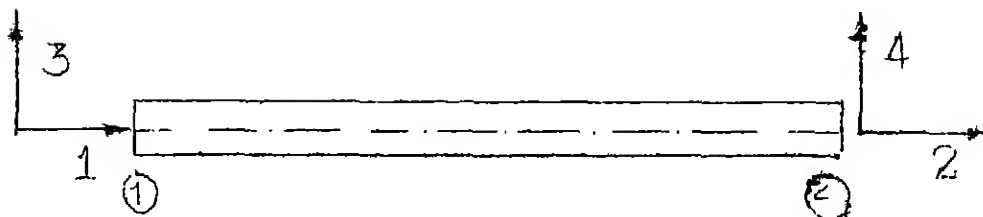
APPENDIX - I

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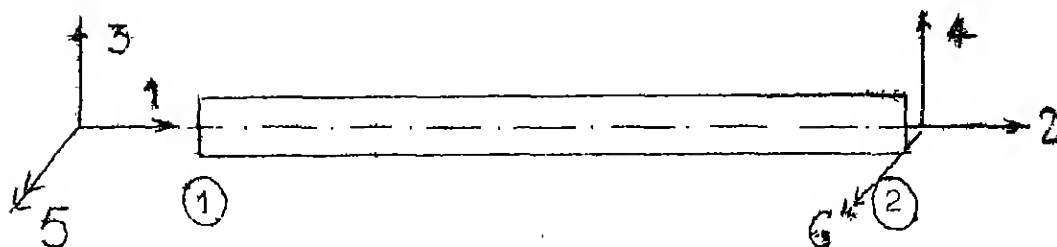
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APPENDIX - II

STIFFNESS AND MASS MATRICES FOR VARIOUS TYPES
OF STRUCTURAL ELEMENTS

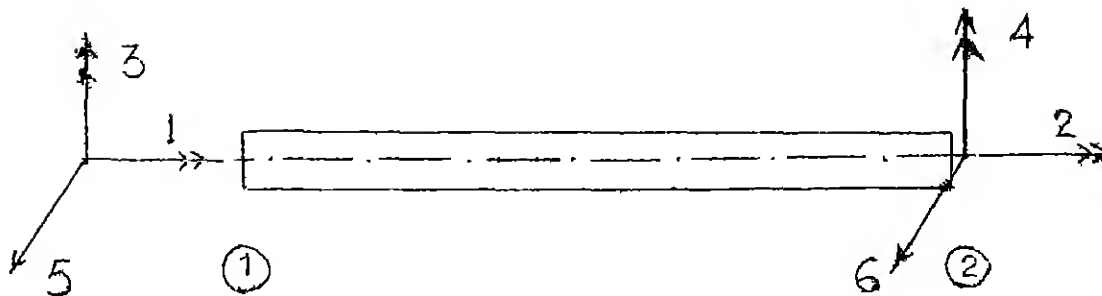
A PLANAR TRUSS ELEMENT

(4. D. O. F)

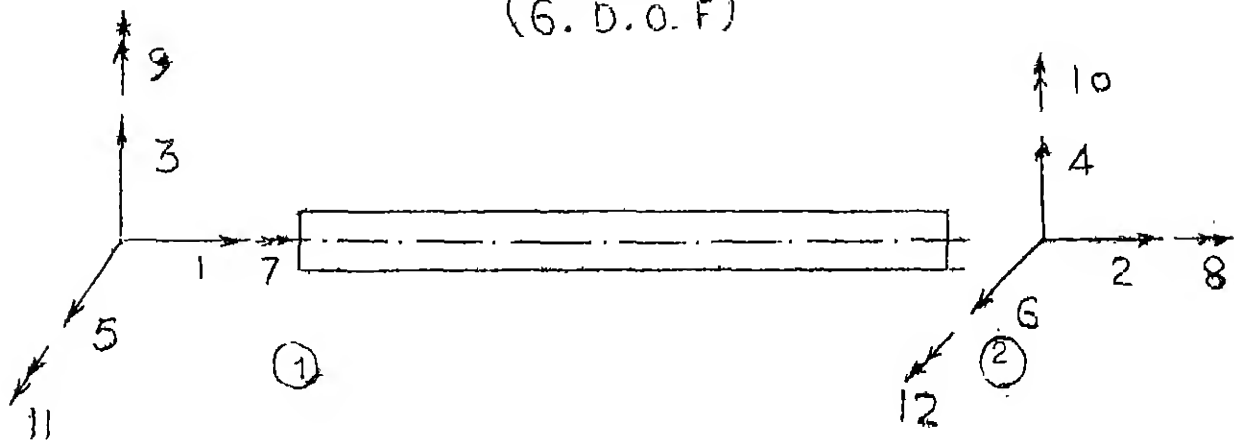
GENERAL BEAM ELEMENT AND A GRID
ELEMENT CONSIDERING INPLANE VIBRATIONS

(6 D. O. F)

STIFFNESS AND MASS MATRICES FOR VARIOUS TYPES OF STRUCTURAL ELEMENTS



A GRID ELEMENT CONSIDERING LATERAL
VIBRATIONS
(6. D.O.F)



A GENERAL SPACE FRAME ELEMENT
(12 D.O.F)

NOTE: DOUBLE ARROWS INDICATE ROTATIONS
AND THE SINGLE ARROW INDICATES
DISPLACEMENTS.

	1	2	3	4
1	$\frac{EA}{L}$	$-\frac{EA}{L}$		
2	$-\frac{EA}{L}$	$\frac{EA}{L}$		
3				
4				

STIFFNESS MATRIX FOR PLANAR TRUSS ELEMENT
(4 D.O.F)

ρ_{AL}

	1	2	3	4
1	$1/3$	$1/6$		
2	$1/6$	$1/3$		
3				
4				

MASS MATRIX FOR PLANAR TRUSS ELEMENT (4 D.O.F.)

	1	2	3	4	5	6
1	$\frac{EA_x}{L}$	$-\frac{EA_x}{L}$				
2	$-\frac{EA_x}{L}$	$\frac{EA_x}{L}$				
3			$\frac{12 EI_z}{L^3}$	$\frac{12 EI_z}{L^3}$	$\frac{6 EI_z}{L^2}$	$\frac{6 EI_z}{L^2}$
4			$-\frac{12 EI_z}{L^3}$	$\frac{12 EI_z}{L^3}$	$-\frac{6 EI_z}{L^2}$	$-\frac{6 EI_z}{L^2}$
5			$\frac{6 EI_z}{L^2}$	$-\frac{6 EI_z}{L^2}$	$\frac{4 EI_z}{L}$	$\frac{2 EI_z}{L}$
6			$\frac{6 EI_z}{L^2}$	$-\frac{6 EI_z}{L^2}$	$\frac{2 EI_z}{L}$	$\frac{AEI_z}{L}$

STIFFNESS MATRIX FOR GENERAL BEAM ELEMENT AND FOR GRID
ELEMENT CONSIDERING INFLANE VIBRATIONS (6 D.O.F)

	1	2	3	4	5	6
1	$1/3$	$1/6$				
2	$1/6$	$1/3$				
3			$13/35$	$9/70$	$(11/210) L$	$- \frac{13}{420} L$
4			$9/70$	$13/35$	$13/420 L$	$- \frac{22}{420} L$
5			$\frac{11}{210} L$	$\frac{13}{420} L$	$L^2/105$	$- \frac{3L^2}{420}$
6			$- \frac{13}{420} L$	$- \frac{22}{420} L$	$- \frac{3L^2}{420}$	$\frac{4}{420} L^2$

MASS MATRIX FOR GENERAL BEAM ELEMENT AND FOR GRID ELEMENT
CONSIDERING INPLANE VIBRATION (6 D.O.F)

1	2	3	4	5	6
GI_x/L	$-GI_x/L$				
$-\frac{GI_x}{L}$	$\frac{GI_x}{L}$				
		$4EI_y/L$	$2EI_y/L$	$-\frac{6EI_y}{L^2}$	$6EI_y/L^2$
		$2EI_y/L$	$4EI_y/L$	$-\frac{6EI_y}{L^2}$	$\frac{6EI_y}{L^2}$
		$-\frac{6EI_y}{L^2}$	$-\frac{6EI_y}{L^2}$	$\frac{12EI_y}{L^3}$	$-\frac{12EI_y}{L^3}$
		$\frac{6EI_y}{L^2}$	$\frac{6EI_y}{L^2}$	$-\frac{12EI_y}{L^3}$	$\frac{12EI_y}{L^3}$

STIFFNESS MATRIX FOR GRID ELEMENT CONSIDERING LATERAL
VIBRATIONS (6 D.O.F)

	1	2	3	4	5	6
1	$\frac{1}{3} I_x / A$	$I_x / 6A$				
2	$I_x / 6A$	$I_x / 3A$				
3			$\frac{4}{420} L^2$	$- L^2 / 140$	$- \frac{11}{210} L$	$- \frac{13}{420} L$
4			$- L^2 / 140$	$\frac{4}{420} L^2$	$\frac{13}{420} L$	$\frac{22}{420} L$
5			$- \frac{11}{210} L$	$\frac{13}{420} L$	$13/35$	$9/70$
6			$\frac{13}{420} L$	$\frac{22}{420} L$	$9/70$	$13/35$

MASS MATRIX FOR GRID ELEMENT CONSIDERING LATERAL VIBRATIONS
(6 D.O.F)

	1	2	3	4	5	6	7	8	9	10	11	12
1	$\frac{EA_x}{L}$	$-EA_x/L$										
2	$-\frac{EA_x}{L}$	EA_x/L										
3			$\frac{12EI_z}{L^3}$	$\frac{12EI_z}{L^3}$							$\frac{6EI_z}{L^2}$	$\frac{6EI_z}{L^2}$
4			$-\frac{12EI_z}{L^3}$	$\frac{12EI_z}{L^3}$							$\frac{6EI_z}{L^2}$	$\frac{6EI_z}{L^2}$
5					$\frac{12EI_y}{L^3}$	$-\frac{12EI_y}{L^3}$			$-\frac{6EI_y}{L^2}$	$-\frac{6EI_y}{L^2}$		
6					$-\frac{12EI_y}{L^3}$	$\frac{12EI_y}{L^3}$			$\frac{6EI_y}{L^2}$	$\frac{6EI_y}{L^2}$		
7							$\frac{GI_x}{L}$	$-\frac{GI_x}{L}$				
8							$-\frac{GI_x}{L}$	$\frac{GI_x}{L}$				
9					$\frac{6EI_y}{L^2}$	$\frac{6EI_y}{L^2}$			$\frac{4EI_y}{L}$	$\frac{2EI_y}{L}$		
10					$-\frac{6EI_y}{L^2}$	$\frac{6EI_y}{L^2}$			$\frac{2EI_y}{L}$	$\frac{4EI_y}{L}$		
11			$\frac{6EI_z}{L^2}$	$-\frac{6EI_z}{L^2}$							$\frac{4EI_z}{L}$	$\frac{2EI_z}{L}$
12			$\frac{6EI_z}{L^2}$	$-\frac{6EI_z}{L^2}$							$\frac{2EI_z}{L}$	$\frac{4EI_z}{L}$

STIFFNESS MATRIX FOR SPACE FRAME ELEMENT (12 D.O.F.)

	1	2	3	4	5	6	7	8	9	10	11	12
1	$1/3$	$1/6$										
2	$1/6$	$1/3$										
3			$13/35$	$9/70$							$\frac{11}{210} L$	$-\frac{13}{420} L$
4			$9/70$	$13/35$							$\frac{13}{420} L$	$-\frac{22}{420} L$
5					$13/35$	$9/70$			$-\frac{11}{210} L$	$\frac{13}{420} L$		
6					$9/70$	$13/35$			$-\frac{13}{420} L$	$\frac{22}{420} L$		
7							$\frac{1}{3} \frac{I_x}{A}$	$I_x/6A$				
8							$I_x/6A$	$I_x/3A$				
9					$-\frac{11}{210} L$	$\frac{13}{420} L$			$\frac{4}{420} L^2$	$-L^2/140$		
10					$\frac{13}{420} L$	$\frac{22}{420} L$			$-\frac{L^2}{140}$	$\frac{4}{420} L^2$		
11			$\frac{11}{210} L$	$\frac{13}{420} L$							$\frac{L^2}{105}$	$-\frac{3L^2}{420}$
12			$-\frac{13}{420} L$	$-\frac{22}{420} L$							$-\frac{3L^2}{420}$	$\frac{4}{420} L^2$

MASS MATRIX FOR A SPACE FRAME ELEMENT (12 D.O.F.)

APPENDIX - III

COMPUTER PROGRAM

This appendix contains a description of each of the Fortran subprograms which makeup the complete program for the free vibration analysis of structures composed of elements described in Appendix - II and the listing of each subprogram is appended at the end.

II-2 The Program:

The complete program for the free vibration analysis of structures consists of (A) a main program which reads the input data, evaluates the lengths of the various members of the structure, generates the starting random vector and prints the various input data. and (B) the other subroutines used are described below:

1 (a) Name of Subroutine: KRMR

(b) Argument List: M, LDF, RDF, NDFS, KL, ML, KR,
MR, TR.

M : Total number of elements

LDF : Local degrees of freedom of an element

RDF : Number of degrees-of-freedom of an element in
system co-ordinates.

NDFS : Number of degrees of freedom of system

KL : Element stiffness matrix in local-coordinate system

ML : Element mass matrix in local-coordinate system

KR : Element stiffness matrix in system co-ordinates

MR : Element mass matrix in system co-ordinates

TR : Transformation matrix for an element

(c) Common Variables: E, A, L, MIX, MIY, MIZ, RO, G, ALPHA,
CX, CY, CZ, PE, QE, NPROB.

E : Modulus of Elasticity of the member

L : Length of the member

MIX : Moment of inertia of the member about x-axis

MIY : Moment of inertia of the member about y-axis

MIZ : Moment of inertia of the member about z-axis

RO : Weight per unit volume of the material of the
member

G : Modulus of rigidity of the member

ALPHA : Angle of inclination of the member

CX, CY and CZ are respectively the X, Y and Z co-ordinates
of the various nodes of the system respectively.

PE and QE : are respectively the left and right hand
ends of a member.

NPROB : is a counter indicating the type of the problem.
 NPROB=1, / indicates that the system consists of truss elements, 2 indicates that the system consists of either the general planar beam element or the grid element for inplane vibrations, 3 indicates that the system consists of grid elements for lateral vibrations and 4 indicates that the system consists of space frame elements.

2 (a) Name of Subroutine: FBMIN

(b) Argument List: M, RDF, NDFS, SR, KR, MR.

Most of them have been defined earlier.

SR : Assembly control table.

(c) Common Variables: SDIFF, AA, BB, NUMER, DENOM, VEC, BX, MIN, MINO, ITER, OMEGA, VALU, VLEO, NVEC.

SDIFF : Scalar; Denotes convergence of Rayleigh quotient function; (MAIN)

AA : Vector; Denotes $[K] \vec{X}$, where $[K]$ is master stiffness matrix and \vec{X} is generalized displacement vector.

BB : Vector; Denotes $[M] \vec{X}$, where $[M]$ is mass matrix and \vec{X} is generalized displacement vector.

NUMER : Scalar; Denotes $\vec{X}^T [K] \vec{X}$.

- DENOM : Scalar; Denotes $\vec{X}^T [M] \vec{X}$.
- VEC : Counter; Denotes the number of eigenvalues--
eigenvectors being searched.
- BX : Matrix ; the ith column of BX represents the
vector $[M] \vec{X}_i$ where $[M]$ is master mass matrix
and \vec{X}_i represents the ith eigenvector.
- MIN : Counter; Denotes the current component number of
the eigenvector which is set to one.
- MINO : Counter; Denotes the old component number of
the eigenvector which was set to one.
- ITER : Counter; Denotes the number of iteration of the
Conjugate gradient method.
- OMEGA : Vector; the components of OMEGA represent the
frequencies (square root of eigenvalues).
- VALU : Vector; the components of VALU represent the
eigenvalues.
- VECO : Matrix, the rows of VECO represent the eigenvectors
- NVEC : Number of eigenvectors used for modal
transformation.

(d) Purpose: To compute the eigenvalues and associated eigenvectors.

3. (a) Name of Subroutine: FANDG

(b) Argument List: for this subroutine the argument list has already been defined.

Purpose: To compute the function (Rayleigh quotient) value and to evaluate the gradient vector to the Rayleigh quotient at any point.

4 (a) Name of Subroutine: PROD

(b) Argument List: Already defined.

Purpose: To compute the product

$$([N][N^T N]^{-1}[N]^T)\vec{W}$$

where $[N] \equiv B\vec{\lambda}$

and \vec{W} is any vector in system co-ordinates,

5(a) Name of Subroutine: INGEN

(b) Argument List: Already defined.

Purpose: To obtain $([N_1]^T [N_1])^{-1}$ from

$$([N_{1-1}]^T [N_{1-1}])^{-1}$$

6. (a) Name of Subroutine : KXKX

(b) Argument List: NDFS, SR, X, KR, MR, LBB

NDFS : Number of degrees of freedom of an element
in system co-ordinates.

SR : Assembly control table

X : Vector in system co-ordinates

KR : Element stiffness matrix in system co-ordinates.

MR : Element mass matrix in system co-ordinates.

LBB : Logical; LBB = .TRUE. provides the normalized
vector $BB \equiv [M] \vec{X}$ (Euclidean norm of the
vector is equal to one).

(c) Common Variables: AA, BB, Y.

AA : Vector; Denotes $[K] \vec{X}$, where $[K]$ is master
stiffness matrix and \vec{X} is generalized displace-
ment vector.

BB : Vector; Denotes $[M] \vec{X}$, where $[M]$ is master
mass matrix and \vec{X} is generalized displacement vector.

Y : Matrix; the *i*th row of the matrix denotes the
displacement vector of the *i*th element corresponding
to any generalized displacement vector \vec{X} in system
co-ordinates.

(d) Purpose: To compute the vectors $[K]\vec{X}$ and $[M]\vec{X}$ where $[K]$ and $[M]$ are respectively the master stiffness and mass matrices of the structure and \vec{X} is any vector in system co-ordinates. Note that master stiffness and mass matrices are not assembled and the desired vectors are obtained by the use of only element stiffness and mass matrices.

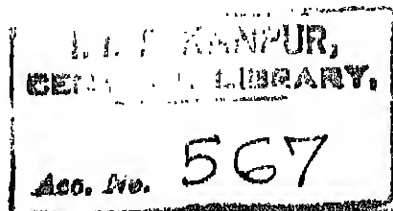
7. (a) Name of Subroutine: TESTS

(b) Argument List: Already defined.

(c) Common Variables: NDFS, NVEC, X, BX, XBX, VEC.

(d) Purpose: Tests for vector space, test for gradient, test for movevector.

A listing of each of the seven fortran subprograms which make up the complete program for the free vibration analysis of various types of structures for IBM 7044 digital computer follows in the subsequent pages.



BFTC MAIN

```

INTEGER PE, QE, RDF, SR
REAL L
REAL MIX, MIY, MIZ
FORMAT(2X, 10E11.6)
2 FORMAT(16F5.1)
FORMAT(8(1X, E8.1))
FORMAT(1X, 25I3)
FORMAT(/, 3X, *% OF MEMBERS = *, I2, 2X, *% OF NODES = *, I3, *LOCAL
* D.O.F = *, I3, 2X, * NO. OF DEGREES OF FREEDOM = *, I3, /, 0, 2X, * NO. OF
* EIGENVALUES AND EIGENVECTORS REQUIRED = *, I3, 2X, * NPROB = *, I3, /)
FORMAT(1H0, 2X, 27H ASSEMBLY CONTROL MATRIX SR, /, 2X, 11H MEMBER NO.,
15X, 19H DEGREES OF FREEDOM, /, 20(6X, I2, 10X, 6(12, 2X), /))
FORMAT(1H0, 2X, 27H ASSEMBLY CONTROL MATRIX SR, /, 2X, 11H MEMBER NO.,
5X, 19H DEGREES OF FREEDOM, /, 20(6X, I2, 10X, 4(12, 2X), /))
FORMAT(1H0, 2X, 27H ASSEMBLY CONTROL MATRIX SR, /, 2X, 11H MEMBER NO.,
5X, 19H DEGREES OF FREEDOM, /, 130(6X, I2, 10X, 12(12, 2X), /))
6 FORMAT(13(1X, F5.1))
9 FORMAT(26F3.1)
3 FORMAT(16F5.1)
DATA PI/3.1412/
COMMON/GMAJN/VECO
COMMON/KAUL/ KR, PD
COMMON/HART/IFF
COMMON/ASINGH/R, RDF, LDF
COMMON/PATIAK/NVEC
COMMON/LEVEL/ E, A, L, MIX, MIY, MIZ, RO, G, ALPHA, CX, CY, CZ, PE, QE, NPROB
DIMENSION E(50), RO(50), CX(50), CY(50), CZ(50), PE(50), QE(50), SR(50, 12) KAUC0280
*, A(50), L(50), VECO(10, 100), MIX(50), MIY(50), MIZ(50), G(50), ALPHA(50) KAUC0290
IFF= 0

```

----- READ INPUT DATA -----

```

READ(5, 5) N, LDF, RDF, NDFS, NVEC, NPROB
READ MODAL CO-ORDINATES
READ(5, 98) (CX(I), I=1, N)
READ(5, 98) (CY(I), I=1, N)
READ(5, 98) (CZ(I), I=1, N)
READ SECTION CHARACTERISTICS
READ(5, 99) (A(I), I=1, 4)
READ(5, 2) (ALPHA(I), I=1, M)
READ(5, 2) (MIX(I), I=1, M)
READ(5, 2) (MIY(I), I=1, M)
READ(5, 2) (MIZ(I), I=1, M)
READ MATERIAL PROPERTIES
READ(5, 2) (E(I), I=1, M)
READ(5, 3) (G(I), I=1, M)
READ(5, 3) (RO(I), I=1, M)

```

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```

KL(2,2)= KL(1,1)
GO TO (10,20,30,40), APPROB
KL(3,3)= (12.0*E(I)*MIZ(I))/(L(I)**3)
KL(3,4)= -KL(3,3)
KL(4,4)= KL(3,3)
KL(3,11)= 0.5*KL(3,3)*L(I)
KL(3,12)= KL(3,11)
KL(4,11)= -KL(3,11)
KL(4,12)= KL(4,11)
KL(5,5)= (12.0*E(I)*MIY(I))/(L(I)**3)
KL(5,6)= -KL(5,5)
KL(5,9)= 0.5*KL(5,6)*L(I)
KL(5,10)= KL(5,9)
KL(6,6)= KL(5,5)
KL(6,9)= 0.5*KL(6,6)*L(I)
KL(6,10)= KL(6,9)
KL(7,7)= G(I)*MIX(I)/L(I)
KL(7,8)= -KL(7,7)
KL(8,8)= KL(7,7)
KL(9,9)= 4.0*E(I)*MIY(I)/L(I)
KL(9,10)= 0.5*KL(9,9)
KL(10,10)= KL(9,9)
KL(11,11)= 4.0*E(I)*MIZ(I)/L(I)
KL(11,12)= 0.5*KL(11,11)
KL(12,12)= KL(11,11)
GO TO 10
KL(3,3)= (12.0*E(I)*MIZ(I))/(L(I)**3)
KL(3,4)= -KL(3,3)
KL(3,5)= 0.5*KL(3,3)*L(I)
KL(3,6)= KL(3,5)
KL(4,4)= KL(3,3)
KL(4,5)= -KL(3,5)
KL(4,6)= -KL(3,5)
KL(5,5)= 4.0*E(I)*MIZ(I)/L(I)
KL(5,6)= 0.5*KL(5,5)
KL(6,6)= KL(5,5)
GO TO 10
KL(1,1)= G(I)*MIX(I)/L(I)
KL(1,2)= -KL(1,1)
KL(2,2)= KL(1,1)
MDI= E(I)*A(I)/L(I)
KL(2,3)= 4.0*MDI
KL(3,4)= 2.0*MDI
KL(3,5)= -6.0*MDI/L(I)
KL(3,6)= 6.0*MDI/L(I)
KL(3,7)= 4.0*MDI
KL(3,8)= KL(3,5)
KL(3,9)= KL(3,6)
KL(3,10)= KL(3,8)
KL(3,11)= KL(3,9)
KL(3,12)= KL(3,11)
KL(4,3)= 4.0*MDI
KL(4,4)= 2.0*MDI
KL(4,5)= -6.0*MDI/L(I)
KL(4,6)= 6.0*MDI/L(I)
KL(4,7)= 4.0*MDI
KL(4,8)= KL(4,5)
KL(4,9)= KL(4,6)
KL(4,10)= KL(4,8)
KL(4,11)= KL(4,9)
KL(4,12)= KL(4,11)
KL(5,3)= 4.0*MDI
KL(5,4)= 2.0*MDI
KL(5,5)= -6.0*MDI/L(I)
KL(5,6)= 6.0*MDI/L(I)
KL(5,7)= 4.0*MDI
KL(5,8)= KL(5,5)
KL(5,9)= KL(5,6)
KL(5,10)= KL(5,8)
KL(5,11)= KL(5,9)
KL(5,12)= KL(5,11)
KL(6,3)= 4.0*MDI
KL(6,4)= 2.0*MDI
KL(6,5)= -6.0*MDI/L(I)
KL(6,6)= 6.0*MDI/L(I)
KL(6,7)= 4.0*MDI
KL(6,8)= KL(6,5)
KL(6,9)= KL(6,6)
KL(6,10)= KL(6,8)
KL(6,11)= KL(6,9)
KL(6,12)= KL(6,11)
KL(7,3)= 4.0*MDI
KL(7,4)= 2.0*MDI
KL(7,5)= -6.0*MDI/L(I)
KL(7,6)= 6.0*MDI/L(I)
KL(7,7)= 4.0*MDI
KL(7,8)= KL(7,5)
KL(7,9)= KL(7,6)
KL(7,10)= KL(7,8)
KL(7,11)= KL(7,9)
KL(7,12)= KL(7,11)
KL(8,3)= 4.0*MDI
KL(8,4)= 2.0*MDI
KL(8,5)= -6.0*MDI/L(I)
KL(8,6)= 6.0*MDI/L(I)
KL(8,7)= 4.0*MDI
KL(8,8)= KL(8,5)
KL(8,9)= KL(8,6)
KL(8,10)= KL(8,8)
KL(8,11)= KL(8,9)
KL(8,12)= KL(8,11)
KL(9,3)= 4.0*MDI
KL(9,4)= 2.0*MDI
KL(9,5)= -6.0*MDI/L(I)
KL(9,6)= 6.0*MDI/L(I)
KL(9,7)= 4.0*MDI
KL(9,8)= KL(9,5)
KL(9,9)= KL(9,6)
KL(9,10)= KL(9,8)
KL(9,11)= KL(9,9)
KL(9,12)= KL(9,11)
KL(10,3)= 4.0*MDI
KL(10,4)= 2.0*MDI
KL(10,5)= -6.0*MDI/L(I)
KL(10,6)= 6.0*MDI/L(I)
KL(10,7)= 4.0*MDI
KL(10,8)= KL(10,5)
KL(10,9)= KL(10,6)
KL(10,10)= KL(10,8)
KL(10,11)= KL(10,9)
KL(10,12)= KL(10,11)
KL(11,3)= 4.0*MDI
KL(11,4)= 2.0*MDI
KL(11,5)= -6.0*MDI/L(I)
KL(11,6)= 6.0*MDI/L(I)
KL(11,7)= 4.0*MDI
KL(11,8)= KL(11,5)
KL(11,9)= KL(11,6)
KL(11,10)= KL(11,8)
KL(11,11)= KL(11,9)
KL(11,12)= KL(11,11)
KL(12,3)= 4.0*MDI
KL(12,4)= 2.0*MDI
KL(12,5)= -6.0*MDI/L(I)
KL(12,6)= 6.0*MDI/L(I)
KL(12,7)= 4.0*MDI
KL(12,8)= KL(12,5)
KL(12,9)= KL(12,6)
KL(12,10)= KL(12,8)
KL(12,11)= KL(12,9)
KL(12,12)= KL(12,11)

```

```

L(6,6)= KL(5,5)
COMPUTATION OF THE MASS MATRIX-----
CONTINUE
D2= DO(I)*A(I)*L(I)
L(1,1)= MD2*1.0/3.0
L(1,1)= MD2*1.0/3.0
L(1,2)= MD2*1.0/6.0
L(2,2)= MD2*1.0/3.0
DO TO (110,120,130,140) , N1208
L(3,3)= MD2*13.0/35.0
L(3,4)= MD2*9.0/70.0
L(3,11)= MD2*L(I)*11.0/210.0
L(3,12)= -MD2*L(I)*13.0/420.0
L(4,4)= MD2*13.0/35.0
L(4,11)= MD2*L(I)*13.0/420.0
L(4,12)= -MD2*L(I)*22.0/420.0
L(5,5)= MD2*13.0/35.0
L(5,6)= MD2*9.0/70.0
L(5,9)= -MD2*L(I)*11.0/210.0
L(5,12)= MD2*L(I)*13.0/420.0
L(6,6)= MD2*13.0/35.0
L(6,9)= -MD2*L(I)*13.0/420.0
L(6,12)= MD2*L(I)*22.0/420.0
L(7,7)= (MD2*MIX(I))/(3.0*A(I))
L(7,8)= 0.5*ML(7,7)
L(8,8)= ML(7,7)
L(9,9)= 4.0*MD2*(L(I)**2)/420.0
L(9,12)= -MD2*(L(I)**2)/140.0
L(10,10)= ML(9,9)
L(11,11)= MD2*(L(I)**2)/1.5.0
L(11,12)= -MD2*3.0*(L(I)**2)/420.0
L(12,12)= ML(9,9)
DO TO 110
L(3,3)= MD2*13.0/35.0
L(3,4)= MD2*9.0/70.0
L(3,3)= MD2*L(I)*11.0/210.0
L(4,4)= ML(3,3)
L(3,5)= -MD2*L(I)*13.0/420.0
L(4,5)= -L(3,6)
L(4,6)= -L(3,5)
L(5,6)= -MD2*(L(I)**2)*9.0/420.0
L(5,5)= MD2*(L(I)**2)/105.0
L(6,6)= MD2*(L(I)**2)*4.0/420.0
DO TO 110
L(1,1)= (MD2*MIX(I))/(3.0*A(I))
L(1,2)= 0.5*ML(1,1)
L(2,2)= ML(1,1)
L(3,3)= MD2*(L(I)**2)*4.0/420.0
L(3,4)= -MD2*(L(I)**2)/140.0
L(3,5)= -MD2*L(I)*11.0/210.0

```

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```

ML(3,6)= -D2*L(1)*13.0/420.0
ML(4,4)= ML(3,3)
ML(4,5)= -ML(3,6)
ML(4,6)= -ML(3,5)
ML(5,5)= MD2*13.0/35.0
ML(5,6)= MD2*9.0/70.0
ML(6,6)= MD2*13.0/35.0
110 CONTINUE
DO 51 K=1,LDF
DO 51 J=K,LDF
KL(J,K)= ML(K,J)
51 ML(J,K)= ML(K,J)
CCC ----- FORMATION OF TRANSFORMATION MATRIX -----
DO 81 I=1,LDF
DO 81 J=1,LDF
81 TR(K,J)= 0.
AA= ALPHA(I)
JT=PE(I)
KT= OE(I)
L1= (CX(KT)-CX(JT))/L(I)
M1= (CY(KT)-CY(JT))/L(I)
N1= (CZ(KT)-CZ(JT))/L(I)
IF (ABS(M1).GT.0.9999) GO TO 90
DENOM= (SQRT(L1**2+M1**2))
L2= (-L1*M1*COS(AA)-M1*SIN(AA))/DENOM
M2= DENOM*COS(AA)
N2= (-M1*M1*COS(AA)+L1*SIN(AA))/DENOM
L3= (L1*M1*SIN(AA)-M1*COS(AA))/DENOM
M3= DENOM*SIN(AA)
N3= (L1*M1*SIN(AA)+L1*COS(AA))/DENOM
GO TO 91
90 L1= 0.0
N1=0.0
L2= -M1*COS(AA)
M2=0.0
N2= SIN(AA)
L3= M1*SIN(AA)
M3=0.0
N3= COS(AA)
91 TR(1,1)= L1
TR(2,1)= L1
TR(3,1)= M2
TR(4,1)= M2
TR(5,1)= M3
TR(6,1)= M3
GO TO (15,25,25,35),IPROB
35 TR(1,3)= M1
TR(1,5)= M1
TR(2,4)= M1

```

```

TR(2,6)= M1
TR(3,1)= L2
TR(3,5)= N2
TR(4,2)= L2
TR(4,6)= N2
TR(5,1)= L3
TR(6,2)= L3
TR(5,3)= L3
TR(6,4)= M3

```

```
DO 92 K=1,6
```

```
KK=K+6
```

```
DO 92 J=1,6
```

```
JJ=J+6
```

```
92 TR(KK,JJ)= TR(K,J)
```

```
GO TO 20
```

```
25 TR(1,3)= M1
```

```
TR(2,4)= M1
```

```
TR(3,1)= L2
```

```
TR(4,2)= L2
```

```
GO TO 80
```

```
15 TR(1,1)= L1
```

```
TR(2,2)= L1
```

```
TR(3,3)= L1
```

```
TR(4,4)= L1
```

```
TR(1,3)= M1
```

```
TR(2,4)= M1
```

```
TR(3,1)= -M1
```

```
TR(4,2)= -M1
```

```
80 CONTINUE
```

```
C TRANSFORM STIFFNESS MATRIX TO REFERENCE COORDINATE SYSTEM
```

```
DO64J=1,LDF
```

```
DO64KK=1,RDF
```

```
SUM=0.0
```

```
DO65KK=1,LDF
```

```
65 SU= SU+KL(J,KK)*TR(KK,K)
```

```
64 TS(J,K)= SUM
```

```
DO67J=1,RDF
```

```
DO67K=1,RDF
```

```
SUM=0.0
```

```
DO68KK=1,LDF
```

```
68 SUM= SU+TR(KK,J)*TS(KK,K)
```

```
67 KQ(J,K)= SUM
```

```
C TRANSFORM MASS MATRIX TO REFERENCE COORDINATE SYSTEM
```

```
66 DO 70 J=1,LDF
```

```
DO70K=1,RDF
```

```
SUM=0.0
```

```
DO71KK=1,LDF
```

```
71 SUM= SUM+L(J,KK)*TR(KK,K)
```

```
70 TS(J,K)= SUM
```

```
69 CONTINUE
```

```

DO 73 K = 1, RDF
  SUM = 0.0
  DO 74 KK = 1, LDF
    SUM = SUM + TR(KK, J) * TS(KK, K)
  74
  73 BR(J, K) = SUM
  72 CONTINUE
  RETURN
END

```

%IBFTC FBMI :

```

SUBROUTINE FBMI (RDF, LDF, SR)
C THIS SUBROUTINE FINDS THE EIGEN VECTORS AND CORRESPONDING EIGEN
C 1-VALUES BY THE MINIMISATION OF THE RAYLEIGH QUOTIENT Z. THE
C 2-CONJUGATE METHOD IS USED FOR MINIMISATION.
  INTEGER RDF, SR, VLC
  REAL DUMPER, DUMPER1, DUMPER2, DUMPER, DUMPER1, DUMPER2, KR, BR
  COMMON /PATHAK/ NVEC
  COMMON /BLOCK/ AA, BB
  COMMON /CBLOCK/ SUMR, DENOM
  COMMON /UBLOCK/ OMEGA
  COMMON /JBLOCK/ JI, JII, ITER
  COMMON /DBLOCK/ VEC
  COMMON /GBLOCK/ DX
  COMMON /GMAIN/ VECO
  COMMON /HARI/ IFF
  DIMENSION SR(50,12), X(100), G(100), S(100), XEN(100), GXEN(100), AA(100)
  *), BR(100), BX(100,10), XBX(100), G(100), DUMPER(100), DUMPER1(100), XEN(100)
  *), G(100), P(100), VALU(100), VECO(10,100), KR(12,12), BR(12,12), OMEGA(100), KAUC
  *TE/P(100), HIT(100), SA(100), CV(100)
  1 FORMAT(1X, 'E15.6')
  13 FORMAT(1H0, 20H RESULTS FROM ITERATION, I8/)
  14 FORMAT(1H0, 15H XEN, GXEN, FEM/(2X, 10E11.6))
  15 FORMAT(1H0, 32H GENERALISED DISPLACEMENT VECTOR,/(2X, 10E11.6))
  16 FORMAT(1H, 10X, 10H1- FUNCTION VALUE, 12X, 9, 5X, 16H1 FUNCTION VALUE
    1, 2X, E20.8)
  17 FORMAT(1H0, 2X, 10H ORTHOGONALITY TEST, E19.8, 5X,
    1 16H GRADIENT LENGTH, E19.8, 5X, 12H MOVE VECTOR, E19.8)
  18 FORMAT(1H0, 12H MOVE VECTOR,/(2X, 10E11.6))
  19 FORMAT(1H0, 40H GENERALISED DISPLACEMENT VECTOR (STARTING POINT),
    1 / (2X, 10E11.6))
  20 FORMAT(1H1, 2X, 32H SEARCH FOR EIGENVECTOR NO. , I2, 2X, 31H AND
    1 CORRESPONDING EIGENVALUE,/(2X, 10E11.6))
  21 FORMAT(1H, 22H TEST FOR VECTOR SPACE)
  22 FORMAT(1H, 18H TEST FOR GRADIENT)
  3 FORMAT(1H, 21H TEST FOR MOVE VECTOR)
  26H CHECK FOR EIGENVALUE NO. , I2,/(2X, 10E11.6))
  32H ***** RESTARTED *****
  34H ***** STEEPEST DESCENT MOVE *****
  3X, 30H FRESH SEARCH FOR EIGENVECTOR NO. , I2, 2X, 31H
  1 SPONDING EIGENVALUE,/(2X, 10E11.6))

```

```

29  FORMAT (///, 1H0, 16H, EIGEN VALUES ARE, /, (2X, E 15.8))
30  FORMAT (///, 1H0, 16H FREQUENCIES ARE, /, (2X, E 15.8))
151  FORMAT(2X,10E11.4)
152  FORMAT(2X,20I5)
    DO 222 I=1,NDFS
    DO 222 J=1,NVEC
222  DX(I,J)=0.0
    DO 333 I=1,NVEC
    DO 333 J=1,NVEC
333  XNM(I,J)=0.0
    SDIFF=1.0E+08
    ITND=2*NDFS
48  CONTINUE
    VEC=1
    WRITE(6,20) VEC
    DO 449 I=1,NDFS
449  X(I)=VFCO(VEC,I)
    XX=0.0
    MIN=0
    MINO=0
    DO 901 I=1,NDFS
    IF ((ABS(X(I))) .LE. XX) GO TO 901
    XX=X(I)
    MIN=I
    MINO=I
901 CONTINUE
    DO 447 I=1,NDFS
447  X(I)=X(I)/XX
    WRITE (6,10) (X(I),I=1,NDFS)
453  IF(VEC.EQ. 1) GO TO 49
    CALL TESTS(NDFS,NVLC,NVLCX,X,DX,XX)
    WRITE (6,21)
    WRITE (6,1) (XBX(I),I=1,JKJ)
49  ITER=0
350 CONTINUE
    IT=1
    NITER=10
    CALL FANDG(NDFS,SR,X,.TRUE.,F1,G)
    NUMF1=NUMBER
    DEMO11=DEFO11
    IF (VEC .EQ. 1) GO TO 85
    CALL TESTS (1,NDFS,NVEC,NVECC,G,X,XBX)
    WRITE (6,22)
    WRITE (6,1) (XBX(I),I=1,JKJ)
4449 IF (VEC .NE. NDFS) GO TO 45
    VALU(VFC)=F1
    OMEGA(VFC)=SQRT(F1)
    DO 67 I=1,NDFS
67  VECO(VEC,I)=X(I)
    WRITE (6,14) (X(I),I=1,NDFS),(G(I),I=1,NDFS),F1
    GO TO 80
85 DO 50 I=1,NDFS
50  S(I)=-G(I)
    ITER=1
    WRITE (6,13) ITER
200 CALL FANDG(NDFS,SR,5,.FALSE.,F,G)

```

NUMBER2=NUMBER

DEL10M2=DEL10M

NUM=0.0

DEN=0.0

DO 51 I=1,NDFS

NUM=NUM+X(I)*AA(I)

51 DEN=DEN+X(I)*BB(I)

U=(NUMBER2*DEN)-(NUM*DEN012)

V=(NUMBER2*DEN041)-(NUMBER1*DEN012)

W=(NUM*DEN011)-(NUMBER1*DEN)

V=V/U

W=W/U

U=1.0

IF(ITER.GT.1.AND.ITER.LT.NITER) GO TO 1200

WRITE (6,1) U,V,W

1200 TS1=-V/(2.0*U)

TS2=((V**2-4.0*U**2)**0.5)/(2.0*U)

TS3=-TS2

ALPHA1=TS1+TS2

ALPHA2=TS1+TS3

IF(ITER.GT.1.AND.ITER.LT.NITER) GO TO 2200

WRITE (6,1) ALPHA1,ALPHA2

2200 IF ((ALPHA1.LT.0.0).AND.(ALPHA2.LT.0.0)) GO TO 76

IF ((ALPHA1.GT.0.0).AND.(ALPHA2.GT.0.0)) GO TO 60

ALPHA=AMAX1(ALPHA1,ALPHA2)

GO TO 76

60 ALPHA=AMIN1(ALPHA1,ALPHA2)

76 IF(ITER.GT.1.AND.ITER.LT.NITER) GO TO 3300

WRITE (6,1) ALPHA

3300 GO TO 71

76 DO 79 I=1,NDFS

79 XEM(I)=X(I)+ALPHA1*S(I)

CALL FANDG (NDFS,SR,XEM,.FALSE.,F1,G)

DO 80 I=1,NDFS

80 XEM(I)=X(I)+ALPHA2*S(I)

CALL FANDG (NDFS,SR,XEM,.FALSE.,F2,G)

IF(ITER.GT.1.AND.ITER.LT.NITER) GO TO 4200

WRITE (6,1) F1,F2

4200 IF (F1.LT.F2) GO TO 77

ALPHA=ALPHA2

GO TO 78

77 ALPHA=ALPHA1

IF(ITER.GT.1.AND.ITER.LT.NITER) GO TO 5200

78 WRITE (6,1) ALPHA

5200 DO 81 I=1,NDFS

81 XEM(I)=X(I)+ALPHA*S(I)

DO 82 I=1,NDFS

82 X(I)=XEM(I)

(6,25)

\$TS(NDFS,NVEC,NVECC,X,3X,XBX)

(6,21)

(6,1) (XBX(I),I=1,J,K)

END

```

      TA=TA+G(I)*G(I)
32  XEM(I)=X(I)+ALPHA*S(I)
      RMAX = 0.0
      IMAX=0
      DO 451 I=1,NDFS
      IF ((ABS(XEM(I))) .LE. RMAX) GO TO 451
      RMAX=ABS(XEM(I))
      IMAX=I
451  CONTINUE
      IF(RMAX .LE. 10.0) GO TO 454
      MINC=MIN
      MIN=IMAX
      DO 452 I=1,NDFS
452  X(I)=XEM(I)
      XX=X(MIN)
      DO 450 I=1,NDFS
450  X(I)=X(I)/XX
      WRITE (6,26)
      GO TO 453
454  CALL FANDG(NDFS,SR,XEM,.TRUE.,FXEM,GXEM)
      NUMER1=NUMER
      DENOM1=DENOM
      IF(ITER.GT.1.AND.ITER.LT.NITER) GO TO 86
      IF(VEC.EQ.1) GO TO 86
      CALL TESTS (NDFS,NVEC,NVECC,XEM,DX,XBX)
      WRITE (6,21)
      WRITE (6,1) (XBX(I),I=1,JKJ)
6665 CALL TESTS (NDFS,NVEC,NVECC,FXEM,DX,XBX)
      WRITE (6,22)
      WRITE (6,1) (XBX(I),I=1,JKJ)
86  GS=0.0
      AG=0.0
      BE=0.0
      DO 53 I=1,NDFS
      GS=GS+S(I)*GXEM(I)
      AG=AG+S(I)*S(I)
53  BE=BE+GXEM(I)*GXEM(I)
      BETA=BE/TA
      AG=SQRT(AG)
      BE=SQRT(BE)
      TEST=GS/(AG*BE)
      IF(ITER.GT.1.AND.ITER.LT.NITER) GO TO 6100
      WRITE (6,17) TEST,BE,AG
6100 DO 54 I=1,NDFS
54  S(I)=-GXEM(I)+BETA*S(I)
      IF(VEC.EQ.1) GO TO 6200
      IF(ITER.GT.1.AND.ITER.LT.NITER) GO TO 8200
      CALL TESTS (NDFS,NVEC,NVECC,S,BX,XBX)
      WRITE (6,23)
      WRITE (6,1) (XBX(I),I=1,JKJ)
8200 IF(ITER.GT.1.AND.ITER.LT.NITER) GO TO 8200
86  WRITE(6,16) F1, FXEM

```

```

200 DIF=F1-FXEM
   IF (DIF .LT. 1.0) GO TO 400
56 DEL=DIF*SDIFF
   DEL=DEL/10.0
   IF (DEL .LT. F1) GO TO 400
   IF ((DEL .LT. F1) .AND. (ABS(DEL) .GT. 1)) GO TO 400
99 DO 55 I=1,NDFS
   G(I)=GXEM(I)
55 X(I)=XEM(I)
   F1=FXEM
   ITER=ITER+1
   IF (ITER .GT. ITNB) GO TO 300
   IF (ITER .GT. NITER) IT=IT+1
   NITER=10*IT
   IF (ITER .GT. 1 .AND. ITER .LT. NITER) GO TO 9200
   WRITE (6,13) ITER
   GO TO 200
300 CONTINUE
   WRITE (6,14) (XEM(I), I=1,NDFS), (GXEM(I), I=1,NDFS), FXEM
   WRITE (6,27) VEC
   GO TO 250
250 CONTINUE
   WRITE (6,14) (XEM(I), I=1,NDFS), (GXEM(I), I=1,NDFS), FXEM
   VALU(VEC)=FXEM
   OMEGA(VEC)=SQRT(FXEM)
   DO 401 I=1,NDFS
401 VECO(VEC,I)=XEM(I)
   CALL KX(X(NDFS,SR,XEM,.FALSE.))
   DO 68 I=1,NDFS
68 CV(I)=AA(I)/BB(I)
   WRITE(6,24) VEC, (CV(J), J=1,NDFS)
240 IF (VEC .EQ. -VEC) GO TO 69
   VEC=VEC+1
   WRITE (6,20) VEC
69 CONTINUE
   CALL KX(X(NDFS,SR,XEM,.TRUE.))
   JJ=VEC-1
   DO 58 I=1,NDFS
58 BX(I,JJ)=CX(I)
   DO 40 I=1,NDFS
40 P(I)=VECO(VEC,I)
28 IF (VEC .GT. 2) GO TO 57
   DEN1=0.0
   DO 64 I=1,NDFS
64 DEN1=DEN1+P(I)*BX(I,1)
   DO 65 I=1,NDFS
65 X(I)=P(I)-DEN1*BX(I,1)
   GO TO 92
57 IF (VEC .GT. 3) GO TO 220
   DEN1=0.0
   DO 59 I=1,NDFS
59 DEN1=DEN1+BX(I,1)*BX(I,2)
   RAS=(1.0-(DEN1)**2)

```

```

X(N(1,1))=1.0/PAS
XNN(1,2)=-DEN1/PAS
XNN(2,1)=-DEN1/PAS
XNN(2,2)=1.0/PAS
GO TO 72
229 VEC=VEC-1
CALL INCEP(NDFS,BX,XNN,DB)
VEC=VEC+1
72 CALL PROD(NDFS,VEC,P,BX,XNN,P,1,P,1,P)
DO 61 I=1,NDFS
61 X(I)=P(I)-C*P(I)
92 XX=0.0
MIN=0
DO 902 I=1,NDFS
IF ((ABS(X(I))) .LE. XX) GO TO 902
XX=X(I)
MIN=I
902 CONTINUE
DO 62 I=1,NDFS
62 X(I)=X(I)/XX
WRITE (6,10) (X(I),I=1,NDFS)
JKU=VEC-1
CALL TESTS (NDFS,NVEC,NVECC,BX,XBX)
WRITE (6,21)
WRITE (6,1) (XBX(I),I=1,JKU)
510 GO TO 49
59 WRITE (6,20) (VALU(I),I=1,NVEC)
WRITE (6,30) (OMEGA(I),I=1,NVEC)
RETURN
END

```

C

\$IBFTC FADG

```

SUBROUTINE FADG(NDFS,SR,Y,LFG,F,G)
INTEGER SR,VEC,DOF
REAL MUER,MNF,NG,MNG,MFG
LOGICAL LFG
CO. ON /SINGH/N,PEF,LDF
CC. ON /PATHAK/NVEC
COM. ON /BLOK/AA,BB
COM. ON /CBLOCK/MUER,DMF
COM. ON /DBLOCK/VEC
CC. ON /CBLOCK/BX
COM. ON /JBLOCK/LIN,ITER
COM. ON /HARI/IFF
DIMENSION X(100),G(100),AA(100),BB(100),C(100,10),NN(100,10),
           Q(10,12),BX(100,10),NMG(100),NG(100),XVC(100)
           J(2X,RE15,B)
           J(18E15,B)

```

KAU06

KAU06


```

      DO 33 I=1,NDFS
33  N(I,K)=BX(I,J)
31  CONTINUE
      VEC=LL+1
      GO TO 11
12  JJ=VEC-1
      DO 26 I=1,NDFS
26  BB(I) = BX(I,JJ)
      CALL INGEN (NDFS,N,NN,BB)
      DO 27 I=1,NDFS
27  N(I,VEC)=BX(I,JJ)
11  CONTINUE
      CALL PROD(NDFS,NVEC,G,N,NN,NG,NNNG,NNNG)
      DO 37 I=1,NDFS
37  G(I)=G(I)-NN*NG(I)
38  G(MIN)=0.0
24  RETURN
      END

```

C. \$IBFTC PROD

```

      SUBROUTINE PROD(NDFS,VEC,G,N,NN,NG,NNNG,NNNG)
C      THIS INTERNAL SUBROUTINE PERFORMS THE THREE MATRIX VECTOR MULTIPLI
C      CATIONS (P(NTN)-1)NT)G
      REAL N,NN,NG,NNNG,NNNG
      INTEGER VEC
      COMMON/DBLOCK/VEC
      DIMENSION G(100),N(100,10),NG(100),NNNG(100),NN(100,10)KAU06
      DO 31 I=1,VEC
      SUM=0.0
      DO 32 J=1,NDFS
32  SUM=SUM+N(J,I)*G(J)
      NG(I)=SUM
31  CONTINUE
      DO 33 I=1,VEC
      SUM=0.0
      DO 34 J=1,VEC
34  SUM=SUM+NN(I,J)*NG(J)
      NNG(I)=SUM
33  CONTINUE
      DO 35 I=1,NDFS
      SUM=0.0
      DO 36 J=1,VEC
36  SUM=SUM+N(I,J)*NNG(J)
      NNNNG(I)=SUM
35  CONTINUE
      RETURN
      END

```

```

C
SIBFTC KXIX
      SUBROUTINE KXIX(NDFS,SR,X,LDX)
C      THIS SUBROUTINE POST MULTIPLIES THE ELEMENT STIFFNESS AND MASS MAT
C      RICES BY THE ELEMENT DISPLACEMENT VECTOR, Y, AND THEN ASSEMBLES THE
C      (K)X AND (M)X VECTORS.
      INTEGER SR,RDF
      REAL X,PR,L
      COMMON/SIBCH/Y,RDF,LDX
C      M IS NUMBER OF ELEMENTS
C      RDF IS NUMBER OF DEGREES OF FREEDOM IN REFERENCE COORDINATE SYSTEM.
C      ND IS NUMBER OF DEGREES OF FREEDOM OF THE SYSTEM
      LOGICAL LDF
      COMMON/UBLOCK/AA,LE
      COMMON/HARI/IFF
      DIMENSION KR(12,12), P(12,12),SR(100,12),X(100),Y(100,12),TS1(100KAU
*,12),TS2(100,12),AA(100),LB(100)
C      ASSIGN SYSTEM DEGREES OF FREEDOM TO ELEMENT DISPLACEMENT VECTOR
      IFF=IFF+1
      REWIND 3
      DO 10 J=1,M
      DO 11 J=1,RDF
        Y(I,J)=0.0
        JJ=SR(I,J)
        IF (JJ) 110,11,11
110    Y(I,J)=X(JJ)
11    CONTINUE
10    CONTINUE
C      CALCULATE (K)X AND (M)X
      DO 12 J=1,M
        IF (IFF.GT.1) GO TO 20
        CALL XDR(KR,J,1)
        WRITE(9) ((KR(J,K),K=1,LDX),J=1,LDX),((P(J,K),K=1,LDX),J=1,LDX)
        GO TO 27
26    READ(3) ((KR(J,K),K=1,LDX),J=1,LDX),((P(J,K),K=1,LDX),J=1,LDX)
27    DO 13 J=1,RDF
      SUM1= 0.0
      SUM2= 0.0
      DO 14 K=1,RDF
        SUM1= SUM1+KR(J,K)*Y(I,K)
14      SUM2= SUM2+P(J,K)*Y(I,K)
13      TS1(I,J)= SUM1
        TS2(I,J)= SUM2
12    CONTINUE
      DO 18 I=1,NDFS
        AA(I)=0.0
18      BB(I)=0.0
      DO 19 K=1,M
      DO 20 KK=1,RDF
        I=SR(K,KK)
        IF (I) 21,20,21
        AA(I)=AA(I)+TS1(K,KK)
        BB(I)=BB(I)+TS2(K,KK)
        CONTINUE
      CONTINUE
      IF (AA(I) .EQ. 0.0) GO TO 22

```

```

23 RAS1=RAS1+BB(I)*DB(I)
   DO 24 I=1,NDFS
24 DB(I)=BB(I)/((RAS1)**0.5)
22 RETURN
   END

```

C
SIBFTC INGEN

```

SUBROUTINE INGEN(NDFS,N,N,X)
C THIS INTERNAL SUBROUTINE GENERATES THE INVERSE OF A Q*Q MATRIX, GIV
C NVEC IS NUMBER OF EIGENVALUES-EIGENVECTORS REQUIRED
  INTEGER VEC
  REAL NN,N,NG,NNNG,NNNNG
  COMMON/PATHAK/NVEC
  COMMON/DBLOCK/VEC
  DIMENSION PN(100),TA(100,10),A1(100,10),A2(100),X(100),NN(100,10),KAUG
  * N(100,10),NG(100),NNNG(100),NNNNG(100)
  CALL PROD(NDFS,NVEC,X,N,N,NG,NNNG,NNNNG)
  DO 10 I=1,NDFS
10 PN(I)=X(I)-NNNNG(I)
   AO=0.0
   DO 11 J=1,NDFS
11 AO=AO+PN(I)*PN(I)
   AO=1.0/AO
   KKK=VEC-1
   DO 12 I=1,KKK
     DO 13 J=1,KKK
13 TM(I,J)=NNNG(I)*NNNG(J)
12 CONTINUE
   DO 14 I=1,KKK
     DO 15 J=1,KKK
15 A1(I,J)=TM(I,J)+AO*TA(I,J)
14 CONTINUE
   DO 16 I=1,KKK
     A2(I)=-AO*TA(I)
   DO 17 I=1,KKK
     DO 17 J=1,KKK
17 NN(I,J)=A1(I,J)
   JJ=VEC
   DO 18 I=1,KKK
     NN(JJ,I)=A2(I)
18 NN(I,JJ)=A2(I)
   NN(JJ,JJ)=AO

```

RETURN
END

C

SUBROUTINE TESTS

SUBROUTINE TESTS(MDEC, VVEC, NVVEC, X, RX, VTX)

INTEGER VEC

COMMON/DR10CV/VEC

DIMENSION X(100), RX(100,10), VBX(100)

CON1=0.0

DO10 I=1, MDEC

10 CON1=CON1+V(I)*V(I)

CON1=((CON1)**0.5)

JJ=VEC-1

DO11 J=1, JJ

SUM=0.0

DO12 I=1, MDEC

12 SUM=SUM+RX(I, J)*V(I)

XRV(J)=SUM

11 CONTINUE

DO13 J=1, JJ

13 XRV(J)=VBX(J)/CON1

RETURN

END

ENDTRY

